

DRAGON Theory for 3-D CANDU Problems

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- 1. Introduction to the Collision Probability Method
- 2. 3–D Collision Probability Calculations
- 3. Solving the Collision Probability Equations
- 4. Condensation and Homogenization Techniques
- 5. Discussion and Conclusion

Introduction to the CP Method

Contents

- The Transport Equation
- The Collision Probability Technique
- Boundary Conditions
- Cross Sections Considerations

The transport equation is a neutron balance equation

$$\mathcal{L}(\vec{r}, E, \vec{\Omega}) = \mathcal{Q}(\vec{r}, E, \vec{\Omega})$$

 \checkmark $\mathcal L$ represents neutron lost from the system:

$$\mathcal{L} = \vec{\Omega} \cdot \vec{\nabla} \Phi(\vec{r}, E, \vec{\Omega}) + \Sigma(\vec{r}, E) \Phi(\vec{r}, E, \vec{\Omega})$$

\checkmark \mathcal{Q} represents neutron created in the system

$$\mathcal{Q}_s = \int dE' d^2 \Omega \Sigma_s(\vec{r}, E' \to E, \vec{\Omega}' \to \vec{\Omega}) \Phi(\vec{r}, E', \vec{\Omega}')$$
$$\mathcal{Q}_f = \chi(\vec{r}, E) \int dE' d^2 \Omega \nu \Sigma_f(\vec{r}, E') \Phi(\vec{r}, E', \vec{\Omega}')$$

Multigroup transport equation

$$\left[\vec{\Omega}\cdot\vec{\nabla} + \Sigma^g(\vec{r}) + D^g(\vec{r})B^2\right]\Phi^g(\vec{r},\vec{\Omega}) = Q_s^g(\vec{r},\vec{\Omega}) + \frac{1}{k}Q_f^g(\vec{r},\vec{\Omega})$$

Scattering source $Q_s^g(\vec{r}, \vec{\Omega})$

$$Q_s^g(\vec{r},\vec{\Omega}) = \sum_{h=1}^G \int d^2 \Omega' \Sigma_s^{hg}(\vec{r},\vec{\Omega}' \to \vec{\Omega}) \Phi^h(\vec{r},\vec{\Omega}')$$

Fission source $Q_f^g(\vec{r}, \vec{\Omega})$

$$Q_f^g(\vec{r}) = \chi^g \sum_{h=1}^G \nu \Sigma_f^h(\vec{r}) \int d^2 \Omega' \Phi^h(\vec{r}, \vec{\Omega}')$$



Transport equation in the absence of external sources is an eigenvalue problem:

- 1. k eigenvalue with imposed leakage ($D^g(\vec{r})B^2$ fixed):
 - Indicates how the fission rate should be modified to make the system critical (reach a non-trivial solution to the transport equation)
- 2. Buckling eigenvalue with imposed k:
 - $D^g(\vec{r})B^2$ represent the amount of leakage required to make the system critical

Boundary conditions are required to close the system

Albedo conditions

$$\phi_{-}(\vec{r}_{S},\vec{\Omega}-2(\vec{N}_{S}\cdot\vec{\Omega}))=\beta(\vec{r}_{S},\vec{\Omega})\phi_{+}(\vec{r}_{S},\vec{\Omega})$$



 $\beta(\vec{r}_S, \vec{\Omega}) = 0$ for void BC $\beta(\vec{r}_S, \vec{\Omega}) = 1$ for reflective BC

COLVE COLLE MONTREAL The Transport Equation



$$\phi_{-}(\vec{r}_{S},\vec{\Omega}) = \phi_{+}(\vec{r}_{S'},\vec{\Omega})$$





Integral transport equation (case without leakage)

■ Flux at a point \vec{r} due to neutrons created at any point $\vec{r}' = \vec{r} - R\vec{\Omega}$ surrounding it

$$\left[-\frac{d}{dR} + \Sigma^g(\vec{r} - R\vec{\Omega})\right] \Phi^g(\vec{r} - R\vec{\Omega}, \vec{\Omega}) = Q^g(\vec{r} - R\vec{\Omega}, \vec{\Omega})$$

Integrate transport equation over R and $\vec{\Omega}$

$$\phi^{g}(\vec{r}) = \int d^{2}\Omega \ e^{-\tau^{g}(R_{S})}(\vec{\Omega} \cdot \vec{N}_{-})\Phi^{g}_{-}(\vec{r}'_{S},\vec{\Omega})\Theta(\vec{r},\vec{r}'_{S},\vec{\Omega})$$
$$+ \int d^{2}\Omega \int_{0}^{R} e^{-\tau^{g}(R')}Q^{g}(\vec{r}',\vec{\Omega})\Theta(\vec{r},\vec{r}',\vec{\Omega})dR'$$

Definitions

$$\begin{split} \phi^g(\vec{r}) &= \int d^2 \Omega \Phi^g(\vec{r},\vec{\Omega}) \\ \tau^g(R) &= \int_O^R \Sigma^g(\vec{r} - R'\vec{\Omega}) dR' \\ \Theta(\vec{r},\vec{r}',\vec{\Omega}) &= \begin{cases} 1 & \text{if } \vec{r} = \vec{r}' + R'\vec{\Omega} \\ 0 & \text{otherwise} \end{cases} \end{split}$$

and $\Phi^g_-(\vec{r}'_S,\vec{\Omega})$ is the incoming angular flux on surface S



Equation for the outgoing flux $\phi_+(\vec{r}_S)$ at S

$$\phi_{+}^{g}(\vec{r}_{S}) = \int d^{2}\Omega(\vec{\Omega}\cdot\vec{N}_{+}) \int_{0}^{R} e^{-\tau^{g}(R')} Q^{g}(\vec{r}',\vec{\Omega}) \Theta(\vec{r}_{S},\vec{r}',\vec{\Omega}) dR'$$
$$+ \int d^{2}\Omega(\vec{\Omega}\cdot\vec{N}_{+})(\vec{\Omega}\cdot\vec{N}_{-}) e^{-\tau^{g}(R_{S})} \Phi_{-}^{g}(\vec{r}'_{S},\vec{\Omega}) \Theta(\vec{r}_{S},\vec{r}'_{S},\vec{\Omega})$$

where

$$\phi^g_+(\vec{r}_S) = \int d^2 \Omega \, \Phi^g_+(\vec{r}_S, \vec{\Omega}) (\vec{\Omega} \cdot \vec{N}_+)$$

- **CP** approximations
 - Solution Divide domain into N_V regions of volume V_i where the cross sections sources are independent of \vec{r} and $\vec{\Omega}$

$$\Sigma^{g}(\vec{r}) = \Sigma_{j}^{g} \qquad \text{for } \vec{r} \in V_{j}$$
$$Q^{g}(\vec{r}, \vec{\Omega}) = \frac{1}{4\pi} q^{g}(\vec{r}) = q_{j}^{g} \qquad \text{for } \vec{r} \in V_{j}$$

(the source is assumed isotropic)

Solution Divide the external boundary S into N_S surfaces of area S_{α} and assume angular flux constant on these surfaces

$$\Phi^g_-(\vec{r}_S,\vec{\Omega}) = \frac{1}{4\pi} \phi^g_{\alpha,-}$$

COLVE COLE MONTREAL The Transport Equation

$$\begin{split} \int d^2 \Omega \int_0^R e^{-\tau^g(R')} Q^g(\vec{r}\,',\vec{\Omega}) \Theta(\vec{r},\vec{r}\,',\vec{\Omega}) dR' = \\ q_j^g \int_{V_i} e^{-\tau^g(R')} \Theta(\vec{r},\vec{r}\,',\vec{\Omega}) \frac{d^3 r}{4\pi |\vec{r}-\vec{r}\,'|^2} \\ \int d^2 \Omega \ e^{-\tau^g(R_S)} \Phi_-^g(\vec{r}\,'_S,\vec{\Omega}) \Theta(\vec{r},\vec{r}\,'_S,\vec{\Omega}) = \\ \phi_{\alpha,-}^g \int_{S_\alpha} e^{-\tau^g(R_S)} \Theta(\vec{r},\vec{r}\,'_S,\vec{\Omega}) \frac{d^2 r}{4\pi |\vec{r}-\vec{r}\,'|^2} \end{split}$$

Transport equations in CP form

$$\phi_i^g = \sum_{\alpha=1}^{N_S} p_{i\alpha}^g \phi_{-,\alpha}^g + \sum_{j=1}^{N_V} p_{ij}^g q_j^g$$
$$\phi_{+,\alpha}^g = \sum_{\beta=1}^{N_S} p_{\alpha\beta}^g \phi_{-,\beta}^g + \sum_{j=1}^{N_V} p_{\alpha j}^g q_j^g$$

where

$$\phi_i^g = \frac{1}{V_i} \int_{V_i} d^3 r \phi^g(\vec{r})$$
$$\phi_{+,\alpha}^g = \frac{1}{S_\alpha} \int_{S_\alpha} d^2 r \phi^g(\vec{r}_S)$$

Four types of probabilities

$$\begin{split} \tilde{p}_{ij}^{g} &= V_{i} p_{ij}^{g} = \int_{V_{i}} \int_{V_{j}} \frac{e^{-\tau^{g}(R)}}{4\pi R^{2}} \Theta_{i} \Theta_{j} d^{3}r' d^{3}r \\ \tilde{p}_{i\alpha}^{g} &= V_{i} p_{i\alpha}^{g} = \int_{V_{i}} \int_{S_{\alpha}} \frac{e^{-\tau^{g}(R_{S})}}{4\pi R_{S}^{2}} (\vec{\Omega} \cdot \vec{N_{-}}) \Theta_{i} \Theta_{\alpha} d^{3}r' d^{2}r \\ \tilde{p}_{\alpha i}^{g} &= \frac{S_{\alpha}}{4} p_{\alpha i}^{g} = \int_{S_{\alpha}} \int_{V_{i}} \frac{e^{-\tau^{g}(R)}}{4\pi R^{2}} (\vec{\Omega} \cdot \vec{N_{+}}) \Theta_{\alpha} \Theta_{i} d^{2}r' d^{3}r \\ \tilde{p}_{\alpha \beta}^{g} &= \frac{S_{\alpha}}{4} p_{\alpha \beta}^{g} = \int_{S_{\alpha}} \int_{S_{\beta}} \frac{e^{-\tau^{g}(R_{S})}}{4\pi R_{S}^{2}} (\vec{\Omega} \cdot \vec{N_{-}}) (\vec{\Omega} \cdot \vec{N_{+}}) \\ &\times \Theta_{\alpha} \Theta_{\beta} d^{2}r d^{2}r' \end{split}$$



Symmetry relations

$$V_i p_{ij}^g = V_j p_{ji}^g$$
$$4 V_i p_{i\alpha}^g = S_\alpha p_{\alpha i}^g$$
$$S_\alpha p_{\alpha\beta}^g = S_\beta p_{\beta\alpha}^g$$

Conservation properties

$$\sum_{\alpha=1}^{N_{\alpha}} p_{i\alpha}^{g} + \sum_{j=1}^{N_{j}} p_{ij}^{g} \Sigma_{j}^{g} = 1$$
$$\sum_{\beta=1}^{N_{\beta}} p_{\alpha\beta}^{g} + \sum_{i=1}^{N_{i}} p_{\alpha i}^{g} \Sigma_{i} = 1$$

Boundary Conditions

Surface flux approximation:

Incoming angular flux on outer surfaces assumed to be independent of $\vec{\Omega}$

Comments

- Outgoing angular flux on outer surfaces integrated over $\vec{\Omega}$
- Angular flux not used at region interfaces
- Approximation for incoming angular flux exact for vacuum BC
- Approximation for incoming angular flux leads to large errors in surface flux when flux is not isotropic



Illustration of approximate BC



Boundary Conditions



Recommendations to reduce errors due to approximate use of BC for a fixed direction

- No special treatment for 2 vacuum BC
- Unfold cell once for 1 vacuum and 1 reflection BC
- Multiply unfold cell for 2 reflection or periodic cell, apply approximate BC on final surfaces and consider results in cell located far from these surfaces.



3)

Example of cell unfolding in direction X



|1|3

3)

Boundary Conditions



Simplifying CP equations using approximate BC Assume

$$\vec{J}_{-}^{g} = \mathbf{A}^{g} \vec{J}_{+}^{g}$$

Final transport equation

$$\vec{\phi}^g = \mathbf{P}^g_{c,vv} \vec{q}^g$$

with the complete collision probability matrix $\mathbf{P}_{c,vv}^g$:

$$\mathbf{P}_{c,vv}^{g} = \left(\mathbf{P}_{vv}^{g} + \mathbf{P}_{vs}^{g}((\mathbf{A}^{g})^{-1} - \mathbf{P}_{ss}^{g})^{-1}\mathbf{P}_{sv}^{g}\right)$$

Two types of multigroup cross-section database can be read by DRAGON

- Mixture macroscopic cross-section
- Isotope microscopic cross-section that contains itself a macroscopic cross-section database

Minimum cross-section requirements for each mixture m

- The multigroup total cross section Σ_m^g
- The isotropic component of the multigroup scattering cross section $\Sigma_{m,s,0}^{h \to g}$ defined as

$$\Sigma_{m,s,0}^{h \to g} = \int_{4\pi} d^2 \Omega^2 \Sigma_{m,s}^{h \to g} (\vec{\Omega}' \to \vec{\Omega}) P_0(\vec{\Omega}' \cdot \vec{\Omega})$$

- The product of the average neutron emitted per fission with the multigroup fission cross section $\nu \Sigma_{m,f}^{g}$
- The multigroup fission spectrum χ_m^g

The linearly isotropic component of the multigroup scattering cross section $\Sigma_{m,s,1}^{h \to g}$

$$\Sigma_{m,s,1}^{h \to g} = \int_{4\pi} d^2 \Omega^2 \Sigma_{m,s}^{h \to g} (\vec{\Omega}' \to \vec{\Omega}) P_1(\vec{\Omega}' \cdot \vec{\Omega})$$

• Required only if B_1 leakage method is used

The transport correction $\Sigma_{m, {\rm tc}}^{g}$

• The transport calculations are performed using transport corrected total ($\tilde{\Sigma}_m^g$) and scattering ($\tilde{\Sigma}_{m,s,0}^{h \to g}$) cross sections

$$\begin{split} \tilde{\Sigma}_{m}^{g} = & \Sigma_{m}^{g} - \Sigma_{m,\text{tc}}^{g} \\ \tilde{\Sigma}_{m,s,0}^{h \to g} = & \Sigma_{m,s,0}^{h \to g} - \delta^{gh} \Sigma_{m,\text{tc}}^{g} \end{split}$$

Takes partially into account the linearly anisotropic scattering contributions

For contribution of multi-neutron production reactions such as

$$\mathbf{X}_{A}^{Z} + \mathbf{n}_{1}^{0} \rightarrow \mathbf{X}_{A-1}^{Z} + 2\mathbf{n}_{1}^{0} + \gamma$$

The scattering cross section must be corrected to take into account this effect

$$\tilde{\Sigma}_{m,s,0}^{h \to g} = \Sigma_{m,s,0}^{h \to g} + 2\delta^{gh} \Sigma_{m,(\mathbf{n},2\mathbf{n})}^{g}$$

where $\Sigma^g_{({\rm n},2{\rm n})}$ is the macroscopic cross section associated with the reaction

Macroscopic cross section data base can be created using

- **•** from the input file using the MAC: module
- from a GOXS file using the MAC: module
- **from a microscopic library using the LIB: module**
- from the homogenization and condensation module EDI:
- from a WIMS-AECL execution using the information available on TAPE16 (side-step method)
- from a HELIOS execution

Microscopic libraries

Many formats can be processed by DRAGON including

- WIMS–AECL format
- MATXS format
- WIMD-D4 format

In DRAGON resonance self-shielding calculations are preformed using the Stamm'ler method



Contents

- Collision Probabilities in 3-D
- Numerical Quadrature and Tracking
- Collision Probability Integration
- Neutron Conservation and CP Normalization



Recall CP approximations

- **Divide domain into** N_V regions of volume V_i
- Assume total cross sections constant inside each region
- Assume sources constant inside each region
 - This has an impact on the selection of the spatial mesh
- Assume sources isotropic inside each region
 This may lead to problem when scattering is highly anisotropic.



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Recall BC approximations

- In the external boundary S into N_S surfaces of area S_α
- Assume flux constant on external surfaces
 - This has an impact on the selection of the spatial mesh
- Assume flux isotropic on external surfaces
 - This may lead to problems when flux is highly anisotropic near external boundaries



Collision Probabilities in 3-D

Collision probability definition

$$\tilde{p}_{ij}^{g} = V_i p_{ij}^{g} = \int_{V_i} \int_{V_j} \frac{e^{-\tau^g(R)}}{4\pi R^2} \Theta_i \Theta_j d^3 r' d^3 r$$

Spherical coordinates for d^3r integral

$$\int_{V_j} d^3 r \Theta_j = \int_{4\pi} d^2 \Omega \int_{R_{i-\frac{1}{2}}}^{R_{i+\frac{1}{2}}} R^2 dR \Theta_j$$

Solution Cartesian coordinates for d^3r' integral

$$\int_{V_i} d^3r' \Theta_i = \int dx' \int dy' \int dR' \Theta_i$$

Collision Probabilities in 3-D

General 3–D geometry for collision probability integration







Final form for collision probability integration

$$\begin{split} \tilde{p}_{ij}^{g} &= \int_{4\pi} \frac{d^{2}\Omega}{4\pi} \int dx' \int dy' \int_{R_{i-\frac{1}{2}}}^{R_{i+\frac{1}{2}}} dR' \int_{R_{j-\frac{1}{2}}}^{R_{j+\frac{1}{2}}} dR \ e^{-\tau^{g}(R',R)} \Theta_{i} \Theta_{j} \\ \tau^{g} &= (R_{i+\frac{1}{2}} - R') \Sigma_{i}^{g} + \sum_{k=i+1}^{j-1} \Delta R_{k} \Sigma_{k}^{g} + (R - R_{j-\frac{1}{2}}) \Sigma_{j}^{g} \end{split}$$

with $\Delta R_k = R_{k+\frac{1}{2}} - R_{k-\frac{1}{2}}$.



Collision Probabilities in 3-D

Notation for optical path







After integration over R' and R, one obtains

$$\begin{split} \tilde{p}_{ij}^g = & \frac{1}{4\pi\Sigma_i^g \Sigma_j^g} \int_0^{4\pi} d^2\Omega \int dx' \int dy' \Theta_i \Theta_j \\ & \times \left[1 - \exp\left(-\tau_{i-\frac{1}{2},i+\frac{1}{2}}^g\right) \right] \exp\left(-\tau_{i+\frac{1}{2},j-\frac{1}{2}}^g\right) \\ & \times \left[1 - \exp\left(-\tau_{j-\frac{1}{2},j+\frac{1}{2}}^g\right) \right] \end{split}$$

with

$$\tau_{i\pm\frac{1}{2},j\pm\frac{1}{2}}^{g} = \Sigma_{i}^{g} \left(R_{i+\frac{1}{2}} - R_{i\pm\frac{1}{2}} \right) + \tau_{i+\frac{1}{2},j-\frac{1}{2}} + \Sigma_{j}^{g} \left(R_{j\pm\frac{1}{2}} - R_{j-\frac{1}{2}} \right)$$


Collision Probabilities in 3-D

Case where $\Sigma_i^g = 0$

$$\tilde{p}_{ij}^g = \frac{1}{4\pi\Sigma_j^g} \int_0^{4\pi} d^2\Omega \int dx' \int dy' \Theta_i \Theta_j$$
$$\times \Delta R_i \exp(-\tau_{i+\frac{1}{2},j-\frac{1}{2}}^g) \left[1 - \exp(-\tau_{j-\frac{1}{2},j+\frac{1}{2}}^g)\right]$$

Case where $\Sigma_i^g = \Sigma_j^g = 0$

$$\tilde{p}_{ij}^g = \frac{1}{4\pi} \int_0^{4\pi} d^2 \Omega \int dx' \int dy' \Theta_i \Theta_j \Delta R_i \Delta R_j \exp(-\tau_{i+\frac{1}{2},j-\frac{1}{2}}^g)$$

Collision Probabilities in 3-D

For \tilde{p}_{ii}^g , one obtains after integration over R' and R

$$\tilde{p}_{ii}^{g} = \frac{1}{2\pi \left(\Sigma_{i}^{g}\right)^{2}} \int_{0}^{4\pi} d^{2}\Omega \int dx' \int dy' \Theta_{i} \Theta_{i}$$
$$\times \left[\tau_{i-\frac{1}{2},i+\frac{1}{2}}^{g} - \left(1 - \exp(-\tau_{i-\frac{1}{2},i+\frac{1}{2}}^{g})\right)\right]$$

For $\Sigma_i^g = 0$ this is simplified to

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$$\tilde{p}_{ii}^g = \frac{1}{4\pi} \int_0^{4\pi} d^2 \Omega \int dx' \int dy' \Theta_i \Theta_i \left(\Delta R_i\right)^2$$

Similar relations are obtained for $\tilde{p}_{i\alpha}^g$ and $\tilde{p}_{\alpha\beta}^g$



3-D Adjuster model in DRAGON



Angular quadrature

- Solution S_n type EQ_N quadrature with $4N_{\Omega}(N_{\Omega}+2)/8$ angular directions $\vec{\Omega}_{1,i}$, $\vec{\Omega}_{2,i}$, $\vec{\Omega}_{3,i}$ and $\vec{\Omega}_{4,i}$
- Global quadrature weight $W_{\Omega} = 2/(N_{\Omega}(N_{\Omega}+2))$
- Tracking in lower half sphere only
- Number of tracking quadrant automatically reduced for symmetric cell
- Select as many angles as possible (Neutrons travel on a straight line)



Cartesian surface quadrature



Cartesian quadrature

- Identify the radius h₊ of the smallest sphere including the geometry
- Select a tracking density ρ_p and define the line spacing δ

$$\delta = \frac{2h_+}{N_p} \qquad N_p = (2\sqrt{\rho_p}h_+) + 1$$

Integration line l_{mn} passes through

$$u_m^x = \left(\frac{2m-1}{2}\right)\delta \qquad u_n^y = \left(\frac{2n-1}{2}\right)\delta$$

Integration weight $W_p = \delta^2$



Comments on Cartesian surface quadrature

In Dragon 3 different planes are selected for each given spatial direction



- Select tracking density as dense as possible
 - Each region must be touched by a maximum number of lines



DRAGON tracking example



DRAGON tracking of a line $l_{m,n}$

- 1. Follow the tracking line as it travels through the cell
- 2. Identify each region i and surface s uniquely
- 3. Identify final region number I associated with set of regions i (a flux region)
- 4. Identify final surface number S associated with sets of surfaces s
- 5. Identify external surfaces and regions i crossed by line
- 6. Compute distance $\tilde{l}_{i,m,n}$ the neutron travels in each region
- 7. Store information on temporary tracking file

Case of symmetric cells (mirror reflection on one side of the cell)

Unfold the cell according to symmetry



Regions in unfolded cells are originally assigned new region numbers

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Angle selection for symmetric cells

- Solution For symmetry with respect to a (y, z) plane, track only in quadrants corresponding to directions $\vec{\Omega}_{1,i}$ and $\vec{\Omega}_{3,i}$
- Solution For symmetry with respect to a (z, x) plane, track only in quadrants corresponding to directions $\vec{\Omega}_{1,i}$ and $\vec{\Omega}_{2,i}$
- Solution For symmetry with respect to a (x, y) plane, track only in all quadrants



Post treatment of tracking file

- Assign to all spatial region *i* its final flux region number combining track segments as required
- For each direction, normalize tracks using

$$l_{i,m,n} = \left(\frac{V_i}{\tilde{V}_i}\right)\tilde{l}_{i,m,n}$$

where

$$\tilde{V}_i = \frac{W_p}{3} \sum_{m=1}^{N_p} \sum_{n=1}^{N_p} \tilde{l}_{i,m,n}$$

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Comments on storage requirements for tracking Maximum number of line segments tracks d_t

$$d_t < 6N(N_\Omega(N_\Omega + 2))\rho_p h_+^2$$

• For $N=1000, h_+=50,
ho_p=20$ t/cm 2 and $N_\Omega=8$ $d_t<20 imes10^9$

tracks segments.



ACR control rods model in DRAGON





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3-D cluster analysis not currently allowed in DRAGON



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For each energy group

- Read a line from tracking file
- Scan this line and add contribution to \tilde{p}_{ii}^g

$$\frac{1}{2} (\Sigma_i^g)^2 \tilde{p}_{ii}^g = \frac{1}{2} (\Sigma_i^g)^2 \tilde{p}_{ii}^g + \sum_n W_n \sum_{m \in i} \left(\tau_{i,n,m}^g - \kappa_{i,n,m}^g \right)$$

where $W_n = W_{\Omega} W_p / 3$ and

$$\begin{split} \kappa^g_{i,n,m} &= \left(1 - \exp\left[-\tau^g_{i,n,m}\right]\right) \\ \tau^g_{i,n,m} &= \Sigma^g_i l_{i,n,m} \quad \text{ with } m \in i \end{split}$$



Scan this line a second time and add contributions to \tilde{p}_{ij}^g

$$\begin{split} \Sigma_i^g \Sigma_j^g \tilde{p}_{ij}^g = & \Sigma_i^g \Sigma_j^g \tilde{p}_{ij}^g \\ &+ \sum_n W_n \sum_{m \in i} \sum_{m' \in j} \kappa_{i,n,m}^g \kappa_{n,m+1,m'-1}^g \kappa_{j,n,m'}^g \end{split}$$

using

$$\kappa_{n,m,m'}^g = \prod_{l=m}^{m'} \exp\left[-\tau_{i,n,l}^g\right]$$



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Finish CP calculations

- Only the contributions with m < m' has been considered
- Symmetrize p_{ij}^g using

$$\tilde{p}_{ij}^g = \tilde{p}_{ij}^g + \tilde{p}_{ji}^g$$



Compute errors on CP conservation rules





CP Normalization



Diagonal Normalization

$$p_{D,ii}^g = p_{ii}^g - \frac{R_i^g}{(\Sigma_i^g)^2 V_i}$$
$$p_{D,\alpha\alpha}^g = p_{\alpha\alpha}^g - \frac{4R_\alpha^g}{S_\alpha}$$

- May result in non-physical negative probabilities
- Cannot be applied to problems involving voided zones



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HELIOS Type Normalization

$$p_{H,ij}^g = (w_i^g + w_j^g) p_{ij}^g \qquad p_{H,\alpha\alpha}^g = (w_\alpha^g + w_\beta^g) p_{\alpha\alpha}^g$$

- Apply conservation laws to above relation
- Solve resulting system for w^g using an iterative process
- Does not lead to negative probabilities and works for void regions
- Default option in DRAGON



Comments on storage requirements for CP matrices

- Number of elements per groups is N^2
- Memory space required for execution is about $5N^2$
- Total disk space required for storage of G group CP is GN^2

Solving the CP Equations

Contents

- The Power Iteration
- The Multigroup Iteration
- Leakage Models

Solving the CP Equations

The multigroup transport equation has the form

$$\vec{\phi} = \mathbf{P}_{c,vv}(\vec{q_s} + \frac{1}{k}\vec{q_f})$$
$$\vec{q_s} = \mathbf{\Sigma}_s \vec{\phi}$$
$$\vec{q_f} = \chi \nu \mathbf{\Sigma}_f \vec{\phi}$$

- $\checkmark \vec{\phi}$ is a $N \times G$ dimensional vector
- **9** $\mathbf{P}_{c,vv}$ is the multigroup CP matrix
 - Diagonal in energy, full in space.
- $\chi \nu \Sigma_f$ is a matrix for neutron production by fission and Σ_s is the scattering matrix
 - Diagonal in space, full in energy





We can decompose the scattering matrix as

$$\Sigma_s = \Sigma_{d,s} + \Sigma_{u,s} + \Sigma_{w,s}$$

with

- $\Sigma_{u,s}$ the up-scattering matrix (lower triangular in energy)
- $\Sigma_{d,s}$ the down-scattering matrix (upper triangular in energy)
- $\Sigma_{w,s}$ the within-group scattering matrix (diagonal in energy)

Solving the CP Equations

Defining W the scattering modified CP matrix

$$\mathbf{W} = (\mathbf{I} - \mathbf{P}_{c,vv} \boldsymbol{\Sigma}_{w,s})^{-1} \mathbf{P}_{c,vv}$$

the transport equation now becomes

$$\vec{\phi} = \mathbf{W}(\Sigma_{d,s}\vec{\phi} + \Sigma_{u,s}\vec{\phi} + \frac{1}{k}\vec{q_f})$$

Assuming

•
$$\vec{q_f}/k$$
 is fixed and $\Sigma_{w,s} = 0$

The above equation can be solved directly from group g = 1 to g = G

Solving the CP Equations

The general equation can be solve using two iteration processes

- The Power iteration illustrated by solving the transport problem with $\Sigma_{w,s} = 0$
- The multigroup iteration illustrated by solving the transport problem with $\vec{q_f}/k$ fixed

The Power Iteration

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- Assume $\Sigma_{u,s} = 0$
- Write an iterative group by broup solution to the transport problem as

$$\vec{\phi}(l) = \mathbf{W}(\mathbf{\Sigma}_{d,s}\vec{\phi}(l) + \frac{\chi}{k(l-1)}\nu\mathbf{\Sigma}_{f}\vec{\phi}(l-1))$$
$$k(l) = \sum_{g=1}^{G}\sum_{i=1}^{N}V_{i}\chi_{i}^{g}\sum_{h=1}^{G}\nu\Sigma_{f,i}^{h}\phi_{i}^{h}(l)$$

with $\vec{\phi}(0)$ a known arbitrary flux distribution and

$$k(0) = \sum_{g=1}^{G} \sum_{i=1}^{N} V_i \chi_i^g \sum_{h=1}^{G} \nu \Sigma_{f,i}^h \phi_i^h(0)$$



The iteration process is repeated until

$$\frac{k(l) - k(l-1)}{k(l)} < \epsilon_1$$

$$\frac{\vec{\phi}(l)}{k(l)} - \frac{\vec{\phi}(l-1)}{k(l-1)} < \epsilon_2 \left| \frac{\vec{\phi}(l)}{k(l)} \right|$$

are both satisfied

- Solve group-by-group this fixed source problem using a Gauss-Seidel strategy

$$\vec{\phi}(l) = \mathbf{W}\left(\mathbf{\Sigma}_{d,s}\vec{\phi}(l) + \mathbf{\Sigma}_{u,s}\vec{\phi}(l-1) + \vec{q}\right)$$

Iterate until

$$\left|\frac{\vec{\phi}^g(l) - \vec{\phi}^g(l-1)}{\vec{\phi}^g(l)}\right| < \epsilon_3$$

Multigroup rebalancing technique

Neutron conservation states that for a converged solution $\phi_i^g(l)$

$$\sum_{i=1}^{N} \Sigma_{i}^{g} V_{i} \phi_{i}^{g}(l) = \sum_{i=1}^{N} R_{i}^{g} V_{i} q_{i}^{g}(l)$$
$$+ \sum_{i=1}^{N} R_{i}^{g} V_{i} \left(\sum_{h=1}^{g} \Sigma_{s}^{h \to g} \phi_{i}^{h}(l) + \sum_{h=g+1}^{G} \Sigma_{s}^{h \to g} \phi_{i}^{h}(l) \right)$$

with $R_i^g = 1 - \sum_{i=1}^N \Sigma_j^g p_{c,ij}^g$

The multigroup problem we effectively solve yields

$$\sum_{i=1}^{N} \sum_{i=1}^{g} V_{i} \phi_{i}^{g}(l) = \sum_{i=1}^{N} R_{i}^{g} V_{i} q_{i}^{g}(l) + \sum_{i=1}^{N} R_{i}^{g} V_{i} \left(\sum_{h=1}^{g} \sum_{s}^{h \to g} \phi_{i}^{h}(l) + \sum_{h=g+1}^{G} \sum_{s}^{h \to g} \phi_{i}^{h}(l-1) \right)$$

To restore conservation at each iteration

Use $\tilde{\phi}_i^g = \alpha^g \phi_i^g$ and assume $\tilde{\phi}_i^g$ satisfies conservation relations, then α^g must satisfy

$$\sum_{h=1}^{G} M^{h \to g} \alpha^{h} = q^{g}$$
$$M^{h \to g} = \sum_{i=1}^{N} R_{i}^{g} V_{i} \left(\sum_{i}^{h} \delta_{gh} - \sum_{s}^{h \to g} \right) \phi_{i}^{h}$$
$$q^{g} = \sum_{i=1}^{N} R_{i}^{g} V_{i} q_{i}^{g}$$

Solve for $\alpha^g(k)$ and rebalance flux



• For appromimate solution $\vec{\Gamma}(l)$

$$\vec{\Gamma}(l) = \mathbf{W}\left(\boldsymbol{\Sigma}_{d,s}\vec{\Gamma}(l) + \boldsymbol{\Sigma}_{u,s}\vec{\Gamma}(l-1) + \vec{q}\right)$$

Define an improved flux distribution for the next iteration using

$$\vec{\phi}(l) = \vec{\Gamma}(l) + \omega(l)\vec{\Delta}(l)$$
$$\vec{\Delta}(l) = \vec{\Gamma}(l) - \vec{\phi}(l-1)$$

and $\omega(l)$ will be computed using a variational procedure

Select $\omega(l)$ in such a way that $\vec{\phi}(l)$ minimizes the tranport functional

$$\mathcal{F}[\vec{\phi}] = \frac{1}{2}\vec{\phi}^T \mathbf{Z}^T \mathbf{Z}\vec{\phi} - \vec{\phi}^T \mathbf{Z}^T \mathbf{W}\vec{q}$$
$$\mathbf{Z} = [\mathbf{I} - \mathbf{W}\left(\mathbf{\Sigma}_{d,s} + \mathbf{\Sigma}_{u,s}\right)]$$

This yield

$$\begin{split} \omega(l) &= -\frac{[\vec{\Delta} - \mathbf{W}\vec{S}_{1}(l)]^{T}[\vec{\Gamma} - \mathbf{W}\vec{S}_{2}(l)]}{[\vec{\Delta} - \mathbf{W}\vec{S}_{1}(l)]^{T}[\vec{\Delta} - \mathbf{W}\vec{S}_{1}(l)]}\\ \vec{S}_{1}(l) &= (\mathbf{\Sigma}_{d,s} + \mathbf{\Sigma}_{u,s})\vec{\Delta}(l)\\ \vec{S}_{2}(l) &= (\vec{q} + \mathbf{\Sigma}_{d,s}\vec{\Gamma}(l) + \mathbf{\Sigma}_{u,s}\vec{\Gamma}(l)) \end{split}$$

Leakage Models

In 3–D, the transport equation can be solved in DRAGON using the B_0 and B_1 leakage models

Both of these models are based on the following factorization of the flux

$$\Phi^g(\vec{r},\vec{\Omega}) \approx \Psi^g(\vec{r},\vec{\Omega}) \exp(i\vec{B}\cdot\vec{r})$$

Transport equation with leakage

$$\begin{split} \vec{\Omega} \cdot \vec{\nabla} \Psi^g(\vec{r},\vec{\Omega}) + [\Sigma^g(\vec{r}) + i\vec{B} \cdot \vec{\Omega}] \Psi^g(\vec{r},\vec{\Omega}) = \\ Q^g_s(\vec{r},\vec{\Omega}) + \frac{1}{k_{e\!f\!f}} Q^g_f(\vec{r}) \end{split}$$

In general we will assume that $k_{e\!f\!f}$ =1


For an infinite homogeneous media the scalar flux and vector current are related to each other according to

$$\vec{\Omega}\Psi^g(\vec{\Omega})=\vec{J}^g(\vec{\Omega})=-iD^g\vec{B}\Psi^g(\vec{\Omega})$$

with D^g is an homogeneous diffusion coefficient

Apply to heterogeneous systems

 $\vec{\Omega} \cdot \vec{\nabla} \Psi^g(\vec{r}, \vec{\Omega}) + [\Sigma^g(\vec{r}) + D^g B^2] \Psi^g(\vec{r}, \vec{\Omega}) = Q_s^g(\vec{r}, \vec{\Omega}) + Q_f^g(\vec{r})$

Find the homogeneous diffusion coefficient compatible with this heterogeneous problem

Leakage Models

- 3
- Assume an heterogeneous solution is known for $B^2 = 0$
- Use this solution to define an equivalent infinite homogeneous problem

$$\Sigma^g \Psi^g(\vec{\Omega}) + i\vec{B} \cdot \vec{J}^g(\vec{\Omega}) = Q_s^g(\vec{\Omega}) + Q_f^g$$

where the cross sections and sources are homogenized using the heterogeneous flux

- Solve the homogeneous problem for D^g and B
- Insert in heterogeneous transport equation and obtain an improved solution.
- Repeat until the iterative procedure is converged

Leakage Models

Solving the homogeneous problem (B_1 model)

Use a 2 terms expansion for the scattering cross section in Legendre polynomials

$$\Sigma^{h \to g}_{s}(\vec{\Omega}' \to \vec{\Omega}) = \Sigma^{h \to g}_{s,0} + 3\Sigma^{h \to g}_{s,1}\vec{\Omega} \cdot \vec{\Omega}'$$

Define

$$\psi^{g} = \int d^{2}\Omega \Psi^{g}(\vec{\Omega})$$
$$\vec{j}^{g} = \int d^{2}\Omega \vec{\Omega}' \Psi^{g}(\vec{\Omega})$$

Leakage Models

- 5
- Insert into the homogeneous transport equation, and integrate to obtain

$$\begin{split} \psi^{g} = &\alpha^{g} \sum_{h} (\Sigma_{s,0}^{h \to g} + \chi^{g} \nu \Sigma_{f}^{h}) \psi^{h} + 3\beta^{g} \sum_{h} \Sigma_{s,1}^{h \to g} \frac{\vec{B} \cdot \vec{j}^{h}}{iB^{2}} \\ \vec{j}^{g} = &\beta^{g} \sum_{h} \left[(\Sigma_{s,0}^{h \to g} + \chi^{g} \nu \Sigma_{f}^{h}) \frac{\vec{B} \psi^{h}}{iB^{2}} + 3\Sigma^{g} \Sigma_{s,1}^{h \to g} \frac{\vec{j}^{h}}{B^{2}} \right] \\ &\alpha^{g} = \frac{1}{B} \arctan\left(\frac{B}{\Sigma^{g}}\right) \\ &\beta^{g} = 1 - \Sigma^{g} \alpha^{g} \end{split}$$

Solve for *B*, ψ^g and \vec{j}^g and compute $D^g = i\vec{B}\cdot\vec{j}^g/B^2\psi^g$





The B_0 homogeneous problem S Assume $\sum_{s,1}^{h \to g} = 0$ and obtain

$$\begin{split} \psi^g = &\alpha^g \sum_h (\Sigma_{s,0}^{h \to g} + \chi^g \nu \Sigma_f^h) \psi^h \\ \vec{j}^g = &\beta^g \frac{\vec{B}}{iB^2} \sum_h (\Sigma_{s,0}^{h \to g} + \chi^g \nu \Sigma_f^h) \psi^h \end{split}$$

The homogeneous diffusion coefficient is then given by

$$D^g = \frac{\beta^g}{\alpha^g}$$

Condensation and Homogenization 1

Contents

- Condensation Technique
- Full Cell Homogenization
- Partial Cell Homogenization and SPH Factors
- Microscopic Cross Section Homogenization

Condensation and Homogenization 2

Condensation and homogenization techniques in DRAGON are based on the following assumptions

Reaction rates are physically meaningfull and should be preserved by the condensation/homogenization procedure

$$R_{i} = \sum_{g} V_{i} \phi_{i}^{g} \Sigma_{i}^{g} = V_{i} \phi_{i} \Sigma_{i}$$
$$R^{g} = \sum_{i} V_{i} \phi_{i}^{g} \Sigma_{i}^{g} = V \phi^{g} \Sigma^{g}$$

The eigenvalue is physically meaningfull and should be preserved by the condensation/homogenization procedure

Condensed transport equation (macrogroup K that includes $g \in G_K$)

$$\vec{\Omega} \cdot \vec{\nabla} \sum_{g \in G_K} \Phi^g(\vec{r}, \vec{\Omega}) + \sum_{g \in G_K} \Sigma^g(\vec{r}) \Phi^g(\vec{r}, \vec{\Omega}) = \sum_{g \in G_K} [Q_s^g(\vec{r}, \vec{\Omega}) + \frac{1}{k} Q_f^g(\vec{r})]$$

Few group version of the same equation is

$$\vec{\Omega} \cdot \vec{\nabla} \Phi^K(\vec{r}, \vec{\Omega}) + \Sigma^K(\vec{r}) \Phi^K(\vec{r}, \vec{\Omega}) = [Q_s^K(\vec{r}, \vec{\Omega}) + \frac{1}{k} Q_f^K(\vec{r}, \vec{\Omega})]$$

It should reproduce condensed multigroup results

The condensation procedure that satisfies our requirements





Multiplying CP transport equation by $\Sigma_i^g V_i$ and summing over all regions *i* yields

$$\sum_{i} \Sigma_i^g V_i \phi_i^g = \sum_{i} V_i [Q_{s,i}^g + \frac{1}{k} Q_{f,i}^g]$$

The equivalent transport equation in a homogeneous infinite cell is

$$\hat{\Sigma}^g V \hat{\phi}^g = V [\hat{Q}^g_{s,i} + \frac{1}{k} \hat{Q}^g_{f,i}]$$

The homogenized and homogeneous transport equations are identical if one selects a flux-volume homogenization technique

Flux-volume homogenization technique





Flux-volume homogenization fails if

The cell is finite (a cell with leakage) and

$$\sum_{j=1}^{N_j} p_{ij}^g \Sigma_j^g \neq 1$$

Partial cell homogenization cell is considered

The heterogeneous N region transport equation homogenized over M regions takes the form

$$\sum_{i \in M_I} V_i \Sigma_i^g \phi_i^g = \sum_{i \in M_I} \sum_J \sum_{j \in M_J} p_{ji}^g (\Sigma^g) [Q_{s,i}^g + \frac{1}{k} Q_{f,i}^g]$$

The ${\cal M}$ region heterogeneous transport equation takes the from

$$V_I \hat{\Sigma}_I^g \hat{\phi}_I^g = \sum_J \hat{p}_{JI}^g (\hat{\Sigma}^g) [Q_{s,J}^g + \frac{1}{k} Q_{f,J}^g]$$

where $\hat{P}_{JI}^{g}(\hat{\Sigma}^{g})$ indicates that the CP are computed using homogenized cross sections

We need

$$\sum_{i \in M_I} V_i \Sigma_i^g \phi_i^g = V_I \Sigma_I^g \phi_I^g$$

and

$$\sum_{J} \hat{p}_{JI}^{g} (\hat{\Sigma}^{g}) [Q_{s,J}^{g} + \frac{1}{k} Q_{f,J}^{g}] = \sum_{i \in M_{I}} \sum_{J} \sum_{j \in M_{J}} p_{ji}^{g} (\Sigma^{g}) [Q_{s,i}^{g} + \frac{1}{k} Q_{f,i}^{g}]$$

to be simultaneously true

The flux-volume homogenization method is not longer adequate because

- There is no simple relation between $\hat{p}_{JI}^g(\hat{\Sigma}^g)$ and $p_{ji}^g(\Sigma^g)$ The alternative here is to use a non-linear process
 - Solution Consider a flux-volume homogenization for ϕ_I^g and Σ_I^g
- Section Redefine the homogeneous flux $\hat{\phi}_I^g$ and cross sections $\hat{\Sigma}_I^g$ as follows

$$\hat{\phi}_I^g = \frac{1}{\mu_I^g} \phi_I^g \qquad \hat{\Sigma}_I^g = \mu_I^g \Sigma_I^g$$

Determine the SPH factors μ_I^g numerically in such a way that

$$\sum_{J} \hat{p}_{JI}^{g} (\hat{\Sigma}^{g}) [Q_{s,J}^{g} + \frac{1}{k} Q_{f,J}^{g}] = \sum_{i \in M_{I}} \sum_{J} \sum_{j \in M_{J}} p_{ji}^{g} (\Sigma^{g}) [Q_{s,i}^{g} + \frac{1}{k} Q_{f,i}^{g}]$$

is true

The definition of the SPH factors automatically ensures

$$\sum_{i \in M_I} V_i \Sigma_i^g \phi_i^g = V_I \Sigma_I^g \phi_I^g = V_I \tilde{\Sigma}_I^g \tilde{\phi}_I^g$$

Microscopic Cross Section

The macroscopic cross section associated with a material is simply the sum over all isotopes of the isotopic macroscopic cross section Σ_I namely

$$\Sigma_i^g = \sum_I \Sigma_{I,i}^g$$

where

$$\Sigma_{I,i}^g = N_{I,i} \sigma_I^g$$

with $N_{I,i}$, the concentration of isotope I in region i

Solution The homogenization and condensation procedure described above remain valid for $\Sigma_{I,i}^g$

Microscopic Cross Section

Since the final concentration of isotope *I* in the cell is given by:

$$N_I = \frac{1}{V} \sum_i N_{I,i} V_i$$

we can define the equivalent homogenized microscopic cross section as:

$$\hat{\sigma}_{I}^{K} = \frac{\mu_{I}^{K}}{N_{I}V\phi^{K}} \sum_{i \in M_{I}} \sum_{g \in G_{K}} N_{I,i}V_{i}\sigma_{I}^{g}\phi_{i}^{g}$$

where the microscopic cross sections now become dependent on the spatial position



Some comments and warning on the CP method

- The sources are assumed constant inside each region
 - Select an adequate spatial discretization This may lead to a large number of region (CP is proportionnal to N²) Some regions may be very small causing problem with tracking
- Select a problem that is not too heterogeneous



Conclusions

- 2
- The angular flux on each external surface are assumed constant and isotropic
 - Try to get rid of external surfaces with re-entrant angular flux
 - Select a model where the region of interest is far from the external surfaces