

The **fixed-point** iterative process and associated error are (where $\rho(\mathbf{P})$ is the spectral radius of \mathbf{P}):

$$\vec{x}^{(0)} = \text{arbitrary}; \quad \vec{x}^{(k+1)} = \mathbf{P}\vec{x}^{(k)} + \vec{b}$$

$$\vec{e}^{(k+1)} = \mathbf{P}\vec{e}^{(k)}; \quad \|\vec{e}^{(k+1)}\| \leq \|\mathbf{P}^{k+1}\| \|\vec{e}^{(0)}\|; \quad \|\mathbf{P}^k\| \approx \rho^k(\mathbf{P})$$

To reduce error by a factor of ϵ , it takes $k \approx \frac{\log(\epsilon)}{\log(\rho(\mathbf{P}))}$ iterations.

Richardson	Jacobi
$(\mathbf{I} - \omega^{(k)}\mathbf{A})\vec{x}^{(k)} + \omega^{(k)}\vec{b}$	$\mathbf{D}^{-1}(\mathbf{D} - \mathbf{A})\vec{x}^{(k)} + \mathbf{D}^{-1}\vec{b}$
GS	SOR
$(\mathbf{D} + \mathbf{L})^{-1}[-\mathbf{U}\vec{x}^{(k)} + \vec{b}]$	$(\mathbf{D} + \omega\mathbf{L})^{-1}\left([(1 - \omega)\mathbf{D} - \omega\mathbf{U}]\vec{x}^{(k)} + \omega\vec{b}\right)$

Table 1: $\vec{x}^{(k+1)}$ for several iterative methods

condition number of a matrix \mathbf{A} is defined as $\kappa(\mathbf{A}) = \|\mathbf{A}\| \|\mathbf{A}^{-1}\|$. If the 2-norm is used, then $\|\mathbf{A}\|_2 = \sigma_1$, $\|\mathbf{A}^{-1}\|_2 = 1/\sigma_m$, and $\kappa_2(\mathbf{A}) = \sigma_1/\sigma_m$; σ_m is the m th singular value of \mathbf{A} .

Let \mathbf{G} be a non-singular **preconditioner**, then $\mathbf{A}\vec{x} = \vec{b}$ can be transformed as $\mathbf{G}^{-1}\mathbf{A}\vec{x} = \mathbf{G}^{-1}\vec{b}$.

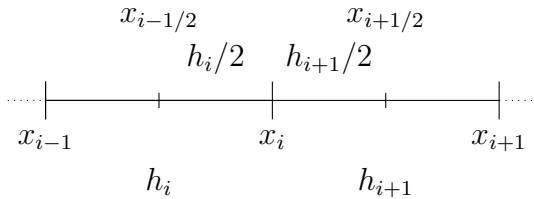
Finite Difference for the DE in 1D we use central difference for the derivative:

$$-\phi_{i-1} + \left(2 + \frac{h^2}{L^2}\right)\phi_i - \phi_{i+1} = h^2 \frac{S_{0,i}}{D} \quad i = 1, 2, \dots, n-1 \quad L \equiv \sqrt{\frac{D}{\Sigma_a}}$$

In 2D we use a 5-point stencil and use central in each dimension for the derivative, giving

$$\nabla^2 \phi_{i,j} = \frac{\phi_{i+1,j} + \phi_{i-1,j} + \phi_{i,j+1} + \phi_{i,j-1} - 4\phi_{i,j}}{h^2}$$

Finite Volume for the DE in 1D uses this grid



The physics values are defined in each cell (cell-centered) and the flux and source are defined at the mesh points (edge-centered).

We integrate the DE in each cell from $x_{i-1/2}$ to $x_{i+1/2}$. At the boundaries we apply the BCs as appropriate to get the $i = 0$ and $i = n$ equations.

The 2D version uses a 2D stencil (not needed for exam). The physics values are again defined in each cell (cell-centered) and the flux and source are defined at the mesh points (edge-centered).

We again integrate over the partial cell, which is now in two dimensions: $x_{i-1/2}$ to $x_{i+1/2}$ and $y_{j-1/2}$ to $y_{j+1/2}$.

We use Gauss Theorem for the streaming term: $-\int_V d\vec{r} [\nabla \cdot (D(\vec{r}) \nabla \phi(\vec{r}))] = \int_S d\vec{S} D(\vec{r}) \frac{\partial}{\partial n} \phi(\vec{r})$.

Critical/eigenvalue equations have the addition of k , allowing us to adjust the equation such that the rhs of the DE becomes $\frac{1}{k} \nu \Sigma_f(x) \phi(x)$.

This means the rhs of our FD or FVM matrix formulation contains flux instead of only a fixed source.

$$A = -\frac{d}{dx} D(x) \frac{d}{dx} + \Sigma_a(x) \quad F = \nu \Sigma_f(x) \quad k = \frac{\text{total production rate}}{\text{total loss rate}}$$

$$\mathbf{A} \vec{\phi}^{(m+1)} = \frac{1}{k^{(m)}} \mathbf{F} \phi^{(m)} \quad k^{(m+1)} = \frac{\int_0^{\bar{a}} F \vec{\phi}^{(m+1)}(x) dx}{\frac{1}{k^{(m)}} \int_0^{\bar{a}} F \vec{\phi}^{(m)}(x) dx}$$

We also add an eigenvalue solver method, like Power Iteration. This uses the idea that $\vec{v}_0 = \gamma_1 \vec{x}_1 + \gamma_2 \vec{x}_2 + \dots + \gamma_n \vec{x}_n$, where \vec{x}_j is the j th eigenvector and γ_j is some constant.

Delayed neutrons let us deal with time variation. β fraction of the total neutrons released from fission are delayed. We use $\rho = \text{reactivity} = \frac{k-1}{k}$, and l is mean neutron lifetime.

$$n(t) = n_0 e^{\frac{(k-1)t}{l}}; \quad T = \text{reactor period} = \frac{l}{k-1}$$

$$\phi(r, t) = v n(t) \psi_1(r); \quad \hat{C}_i(r, t) = C_i(t) \psi_1(r)$$

where ψ_1 is the fundamental mode solution of $\nabla^2 \psi_n + B_g^2 \psi_n = 0$

PRKE

$$\frac{dn(t)}{dt} = \frac{\rho(t) - \beta}{\Lambda} n(t) + \sum_{i=1}^6 \lambda_i C_i(t); \quad \frac{dC_i(t)}{dt} = \frac{\beta_i}{\Lambda} n(t) - \lambda_i C_i(t)$$

$C_i(t)$ = delayed neutron concentration from i th precursor; Λ = effective n lifetime = $(\nu \nu \Sigma_F)^{-1}$.

Runge Kutta methods use stages to avoid higher order derivatives to get better accuracy. Two stage:

$$U^{n+1/2} = U^n + \frac{1}{2} k f(U^n) \quad U^{n+1} = U^n + k f(U^{n+1/2}) \quad U^{n+1} = U^n + k f(U^n + \frac{1}{2} k f(U^n))$$

Monte Carlo uses random numbers, sampling rules, tallies, and statistics

$$\mu = E(x) = \int x f(x) dx; \quad \bar{x} = \frac{1}{N} \sum_{i=1}^N x_i \quad \lim_{N \rightarrow \infty} \bar{x} \rightarrow \mu$$

$$\sigma^2 = E(x^2) - \mu^2; \quad S_x^2 \approx \bar{x}^2 - \bar{x}^2; \quad S_{\bar{x}}^2 = \frac{S_x^2}{N}$$

$$\begin{aligned}
p\{a \leq x \leq b\} &= \int_a^b p(x)dx ; & p(x) &\geq 0 ; & \int_{-\infty}^{\infty} p(x)dx &= 1 \\
p(x = x_k) &= p_k \equiv p(x_k) , k = 1, \dots, N ; & p_k &\geq 0 ; & \sum_{k=1}^N p_k &= 1 \\
P\{x' \leq x\} &= P(x) = \int_{-\infty}^x p(x')dx' ; & P(-\infty) &= 0 , & P(\infty) &= 1 \\
P\{x' \leq x\} &= P_k \equiv P(x_k) = \sum_{j=1}^k p_j , k = 1, \dots, N ; & P_0 &= 0 , & P_N &= 1
\end{aligned}$$

Deal with normalization using $p(x) = \frac{g(x)}{G(\infty)}$, where $p(x)$ is the numerical PDF, $g(x)$ is the physical PDF, and $G(x)$ is the physical CDF.

With rejection sampling, create $g(x) \geq p(x)$ for all x ; Generate pair of random variables, (ξ, η) ; $x' = G^{-1}(\xi)$; If $\eta < p(x')/g(x')$, accept x' ; Else, reject x' .

Probability of distance to collision

$$\begin{aligned}
p_c(s) &= \Sigma_t(s) e^{-\Sigma_t(s)s} ds \\
P_c(n) &= \int_0^s \Sigma_t(s) e^{-\Sigma_t(s)s'} ds' = -e^{-\Sigma_t(s)s'} \Big|_0^s = 1 - e^{-\Sigma_t(s)s}
\end{aligned}$$

After a collision, which nuclide from $p_i = \frac{\Sigma_{t,i}}{\Sigma_t}$ and which reaction from $p_x = \frac{\Sigma_{x,i}}{\Sigma_{t,i}}$.

In scattering, the new direction is $(\sin(\phi) \cos(\theta), \sin(\phi) \sin(\theta), \cos(\theta)) = (\sqrt{1 - \mu^2} \cos(\theta), \sqrt{1 - \mu^2} \sin(\theta), \mu)$.