



Time-Domain Self-Force Computations Using Pseudospectral Methods

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Motivation

The metric perturbations generated by a point particle moving in a Schwarzschild black hole

$$\square_* \mathbf{U} + \mathbb{A} \partial_t \mathbf{U} + \mathbb{B} \partial_{r^*} \mathbf{U} + \mathbb{C} \mathbf{U} = \mathbf{F} \delta[r - r_p(t)],$$

Scalar charged particle falling in a geodesic of a Schwarzschild MBH spacetime.

$$\square \Phi = g^{\mu\nu} \nabla_\mu \nabla_\nu \Phi = -\rho 4\pi$$

$$\rho = -4\pi q \int \delta_4(x - z(\tau)) d\tau$$

Problems: Distributional source term and divergence of the field at the particle location

The retarded field can be decomposed into spherical harmonics:

$$\Phi^{ret} = \sum_{l=0}^{\infty} \sum_{m=-l}^l \Phi^{lm}(t, r) Y^{lm}(\theta, \varphi)$$

The equation for each harmonic coefficient:

$$(-\partial_t^2 + \partial_{r^*}^2 - V_l)\psi^{lm} - S^{lm} = 0$$

$$V_l = \left(1 - \frac{2M}{r}\right) \left[\frac{2M}{r^3} + \frac{l(l+1)}{r^2}\right] \quad \Phi^{lm} = \frac{\psi^{lm}}{r}$$

$$\rho \rightarrow S^{lm} = A^{lm} \delta[r^* - r_p^*(t)]$$

The Particle-without-Particle Scheme

The mode-sum regularization scheme provides an analytic expression for the field singularities.

$$\begin{array}{c} \boxed{\text{Computed Analytically}} \\ \downarrow \\ \Phi^{ret} = \Phi^S + \Phi^R \quad \left\{ \begin{array}{l} \square \Phi^S = -4\pi q \delta(z) \\ \square \Phi^R = 0 \end{array} \right. \\ \uparrow \\ \boxed{\text{Computed Numerically}} \end{array}$$

$$\mathcal{F}_\alpha = q(\nabla_\alpha \Phi^{ret} - \nabla_\alpha \Phi^S) = q\nabla_\alpha \Phi^R$$

To compute the Φ^{ret} using PSC methods we have developed a scheme that removes the singularity associated with the particle

$$\mathcal{U} \Rightarrow \Phi^{ret}$$

Numerical techniques to compute the field modes

Frequency domain

- *Fourier harmonic decomposition:*

Solve ODEs

- *Sum over the Fourier harmonics:*

Difficulties for handling high eccentric orbits

Time domain

- *Solve the PDE for each field mode*
- *Handles in the same way circular and eccentric orbits*
- *Numerically expensive due to the large scale variance in the solutions*

Three approaches for solving PDEs in time-domain

- *Finite differences*
- *Finite elements (FE)* → *Easiest to code but are computationally expensive*
- *Spectral methods (SM)* → *Both expand the solution in basis functions and use multidomain grids*

- *FE suited to irregular geometries.*
- *SM use fewer subdomains than FE and for sufficiently regular domains, are generally faster and/or more accurate.*

[see e.g. H. P. Pfeiffer et al. 2003]

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Outlook

- *Basics about pseudospectral methods*
- *Polynomial expansion and discretization points*
- *How to apply pseudospectral-collocation methods to the EMRI problem: Scalar case in Schwarzschild space-time*
- *Examples: Circular & eccentric case*
- *Results*

Consider an arbitrary system of hyperbolic partial differential equations (PDEs) defined on $\Omega \subset \mathbb{R}^d$

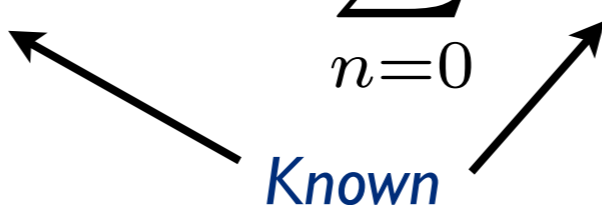
$$L[\mathcal{U}](x) = S(x) \quad x \in \Omega$$

with boundary conditions

$$H[\mathcal{U}](x) = 0 \quad x \in \partial\Omega$$

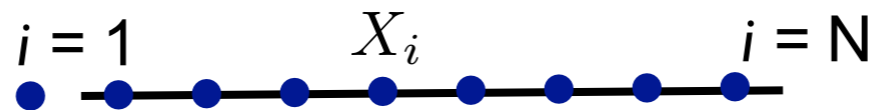
The spectral representation

In 1D spatial dimension pseudospectral collocation methods (PSCM) are based on expansions of every evolved field $\mathcal{U}(x)$ in terms of suitable basis functions $T_n(X)$ with (spectral) coefficients a_n

$$\mathcal{U}(X) \equiv \mathcal{U}_N(X) = \sum_{n=0}^N a_n T_n(X)$$


Known

We can derive a_n from the values of $T_n(X)$ at the (discretization) collocation points $X = X_i$ ($i = 1 \dots N$)



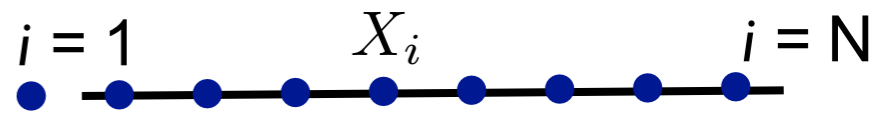
$$\mathcal{U}(X_i) = \sum_{n=0}^N a_n T_n(X_i) \longrightarrow a_n = \frac{[\mathcal{U}_n(X_i), T_n(X_i)]}{|T_n(X)|}$$

$$\mathbb{M}_{ij} \mathcal{U}(X_i) = a_j$$

The physical representation

$\mathcal{U}(x)$ is approximated employing the Lagrange Cardinal functions

$\mathcal{C}_i(x_j) = \delta_{ij}$, associated with $\{T_n(X)\}$.

$$\mathcal{U}_N(X) = \sum_{n=0}^N \mathcal{U}(X_i) \mathcal{C}_i(X)$$


Change between representations employing matrix multiplication transformation

$$\mathcal{U}(X_i) = \mathbb{M}_{ij}^{-1} a_j$$

The derivatives of $T_n(X)$ and $C_i(X)$ are known analytically

$$\partial_X^{(m)} \mathcal{U}_N(X) = \sum_{n=0}^N a_n \partial_X^{(m)} T_n(X)$$

$$\partial_X^{(m)} \mathcal{U}_N(X_i) = \sum_{i=0}^N \sum_{j=0}^N \partial_X^{(m)} C_i(X_j) \mathcal{U}_N(X_j) C_i(X)$$

The derivatives are obtained through a matrix multiplication

Interpolation Error

The error in interpolating the solution is given by Cauchy interpolation error

$$u - u_N(X) = \frac{1}{N+1!} u^{N+1}(\xi) \prod_{i=0}^N (X - X_i)$$

controlled by changing the location of the collocation points,

$$\text{Error} = \log_{10} |a_N|$$

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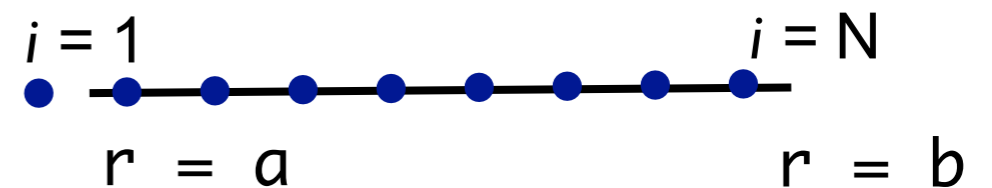
$$\text{Error} = \log_{10} |a_N|$$

The solutions converge exponentially with N

Time step condition

The domain of dependence of the system of hyperbolic equations evolve at finite velocity which leads to causality restrictions: Courant-Friedrichs-Lax conditions

$$\Delta t_{CFL} \sim \frac{\pi^2 |b - a|}{4N^2}$$

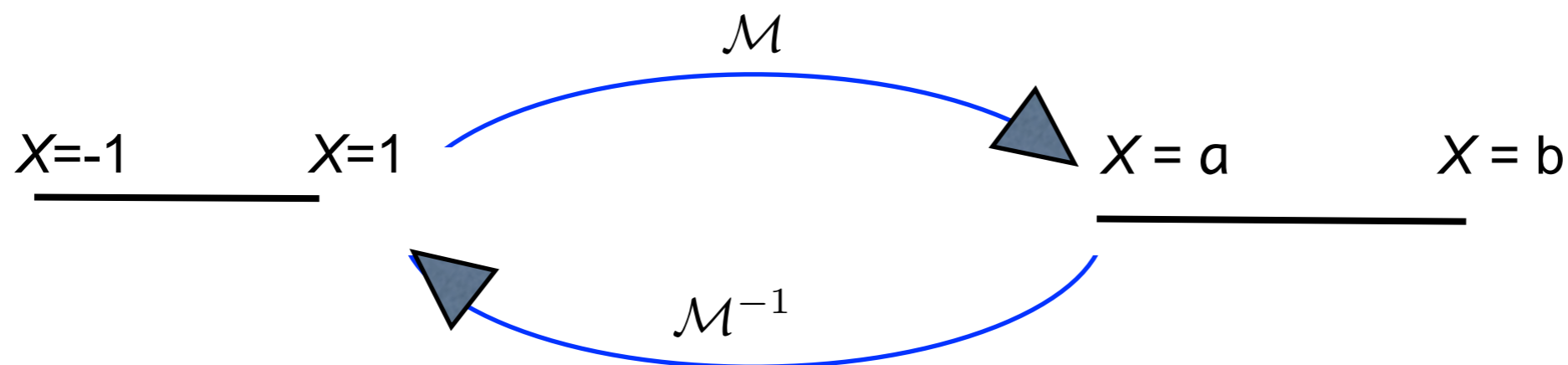


Expansion Basis & and discretization points

The solutions can be discretized using an expansion in a basis of Chebyshev polynomials

$$T_n(X) = \cos \left(n \cos^{-1}(X) \right), \quad X \in [-1, 1].$$

The domain of definition of $T_n(X)$ can always be mapped to the spatial (sub)domain of our problem



Chebyshev series can be expressed as a Fourier cosine series

$$\begin{aligned} X : \quad [0, 2\pi] &\longrightarrow [-1, 1] \\ \theta &\longrightarrow X(\theta) = \cos(\theta) \end{aligned}$$

$$\mathcal{U}(X) = \sum_{n=0}^{\infty} a_n T_n(x) \Leftrightarrow \mathcal{U}(\cos \theta) = \sum_{n=0}^{\infty} a_n \cos(n\theta)$$

The Fourier series of $\mathcal{U}(\cos \theta)$ have exponential convergence, unless $\mathcal{U}(X)$ is singular.

Matrix multiplications can be performed using a FFT algorithm:

$\mathcal{O}(N \ln(N))$ operations instead of $N \times N$ operations needed in a direct matrix multiplication

$$\partial_{r^*} : \{\mathbf{U}_i\} \xrightarrow{FFT} \{\mathbf{a}_n\} \xrightarrow{\partial_{r^*}} \{\mathbf{b}_n\} \xrightarrow{FFT} \{(\partial_{r^*} \mathbf{U})_i\}$$

The set of collocation points that minimises the Cauchy interpolation error corresponds to the zeros of the Chebyshev polynomial or alternatively to the extrema of its derivative: Lobatto-Chebyshev grid

$$(1 - X^2)T'_N(X) = 0$$

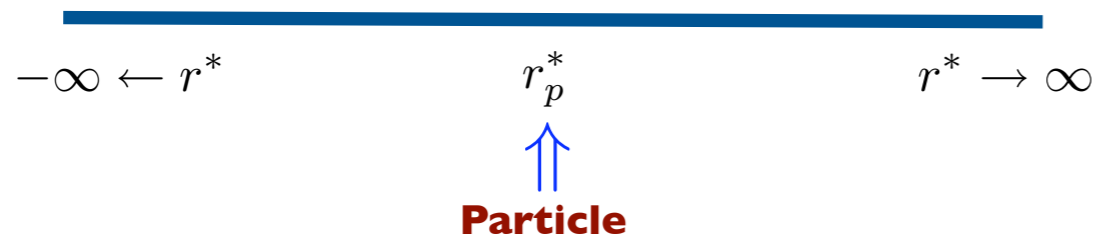
$$X_i = -\cos\left(\frac{\pi i}{N}\right) \quad (i = 0, 1, \dots, N)$$

PSCM & The EMRI problem

The Particle-without-Particle Scheme

$$(-\partial_t^2 + \partial_{r^*}^2 - V_\ell)\psi^{\ell m} = A^{\ell m} \delta[r^* - r_p^*(t)]$$

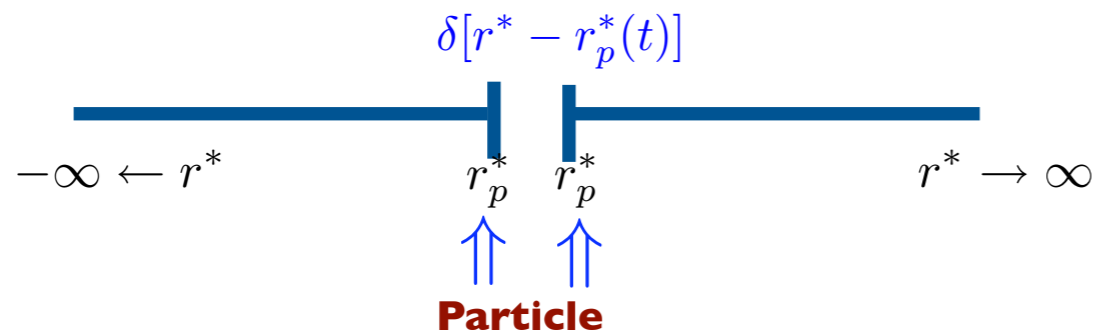
$$\delta[r^* - r_p^*(t)]$$



The computational domain is split and the particle is set at the interface between 2 subdomains

The Particle-without-Particle Scheme

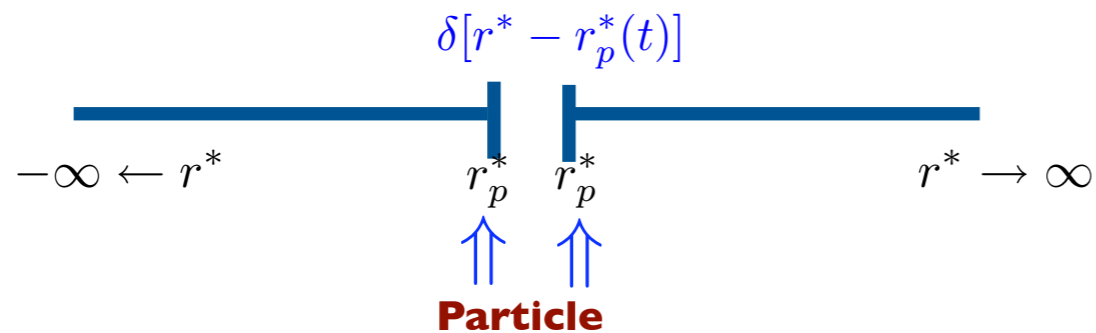
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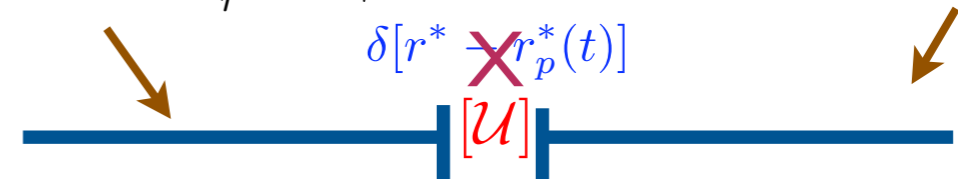
$$U = U_- \Theta(r_p^* - r^*) + U_+ \Theta(r_p^* - r^*)$$

$$\partial_t U_\pm = \mathbb{A} \cdot \partial_{r^*} U_\pm + \mathbb{B} \cdot U_\pm$$

+

$$[U] = \lim_{r^* \rightarrow r_p^*} U_+ - \lim_{r^* \rightarrow r_p^*} U_-$$

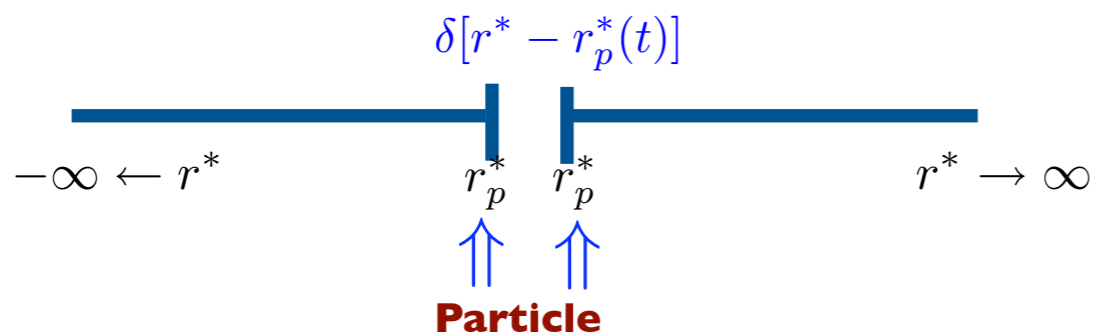
$$\partial_t U_- = \mathbb{A} \cdot \partial_{r^*} U_- + \mathbb{B} \cdot U_- \qquad \partial_t U_+ = \mathbb{A} \cdot \partial_{r^*} U_+ + \mathbb{B} \cdot U_+$$



The PwP technique ensures smooth solutions

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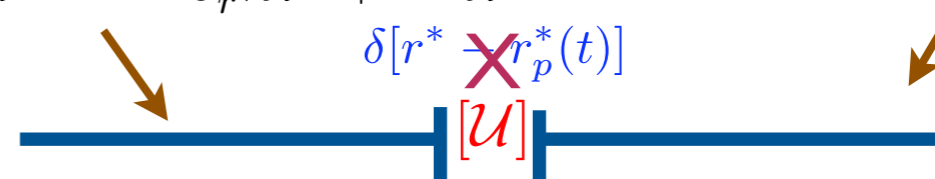
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The PwP technique ensures smooth solutions

The discontinuities on hyperbolic equations propagate along the characteristics.

$$U = (\psi^{\ell m}, \phi^{\ell m}, \varphi^{\ell m})$$

$$\partial_t U = \mathbb{A} \cdot \partial_{r^*} U + \mathbb{B} \cdot U$$

$$\mathbb{A} = \begin{pmatrix} 0 & 0 & 0 \\ 0 & 0 & 1 \\ 0 & 1 & 0 \end{pmatrix} \quad \mathbb{B} = \begin{pmatrix} 0 & 1 & 0 \\ -V_\ell & 0 & 0 \\ 0 & 0 & 0 \end{pmatrix}$$

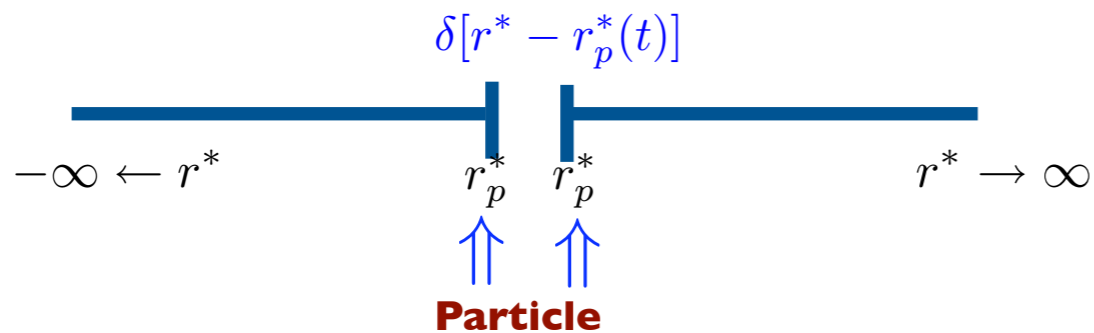
$$\psi^{\ell m} = r \Phi^{\ell m}$$

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The Particle-without-Particle Scheme

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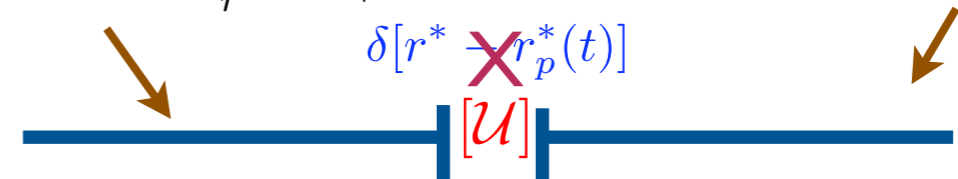
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$$\begin{aligned} \psi^{\ell m} &= r \Phi^{\ell m} \\ \phi^{\ell m} &= \partial_t \psi^{\ell m} \\ \varphi^{\ell m} &= \partial_{r^*} \psi^{\ell m} \end{aligned}$$

The jumps are enforced employing two different methods:

1. The penalty method.
2. Communication of the characteristic fields

The Particle-without-Particle Scheme

1. The penalty method:

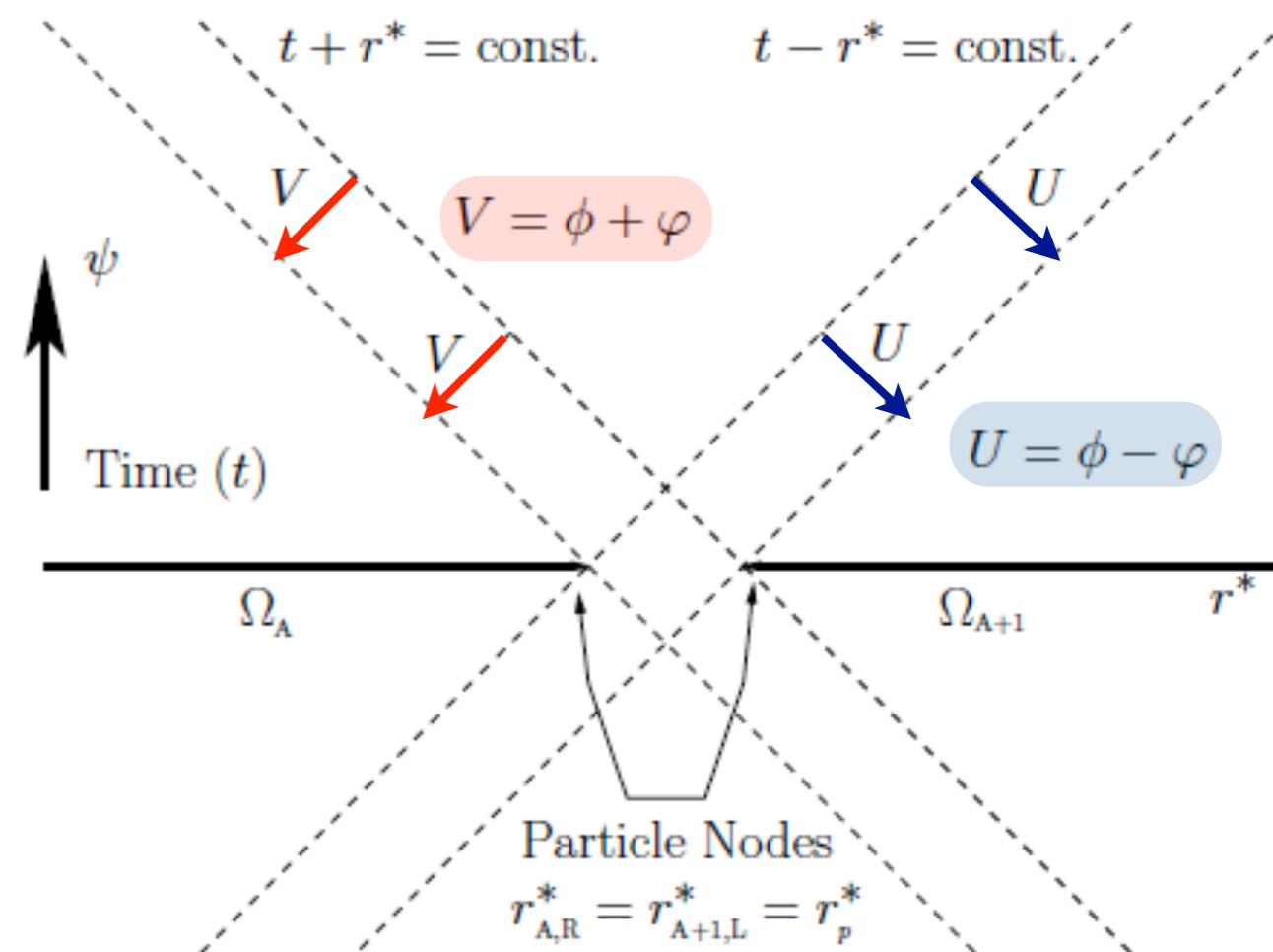
The system is dynamically driven to fulfil a set of additional conditions.

$$\partial_t \mathcal{U}_{\pm} = \mathbb{A} \cdot \partial_{r^*} \mathcal{U}_{\pm} + \mathbb{B} \cdot \mathcal{U}_{\pm} + \eta(\tau_u[\mathcal{U}])$$

2. The direct communication of the characteristic fields:

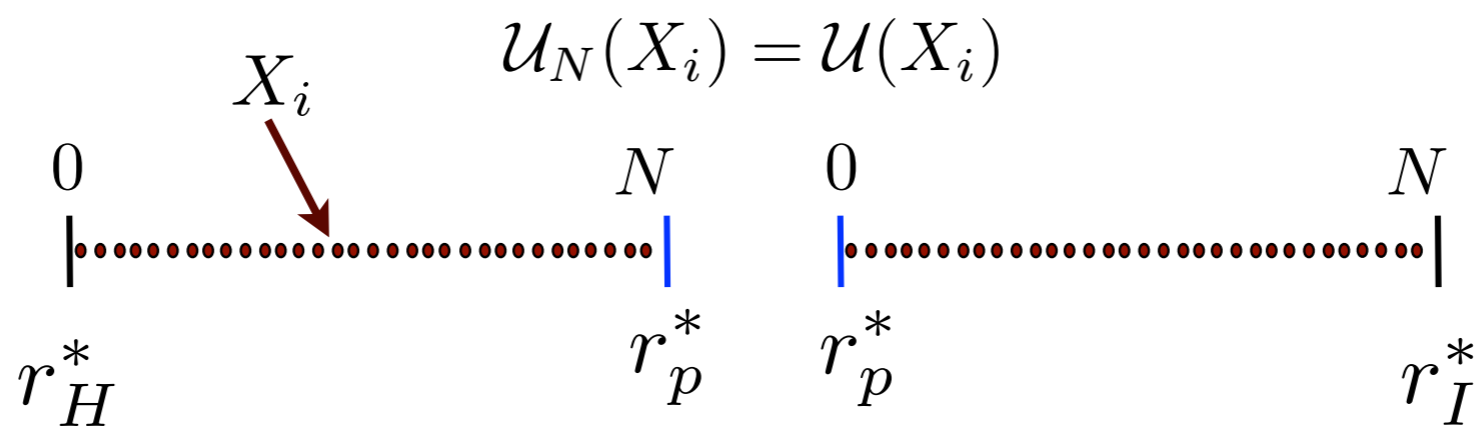
We pass the value of the characteristic fields.

$$\begin{aligned} \psi^{lm} &= r \Phi^{lm} \\ U^{lm} &= \phi^{lm} - \varphi^{lm} \\ V^{lm} &= \phi^{lm} - \varphi^{lm} \\ &\downarrow \\ \mathcal{U} &= (\psi^{lm}, U^{lm}, V^{lm}) \end{aligned}$$



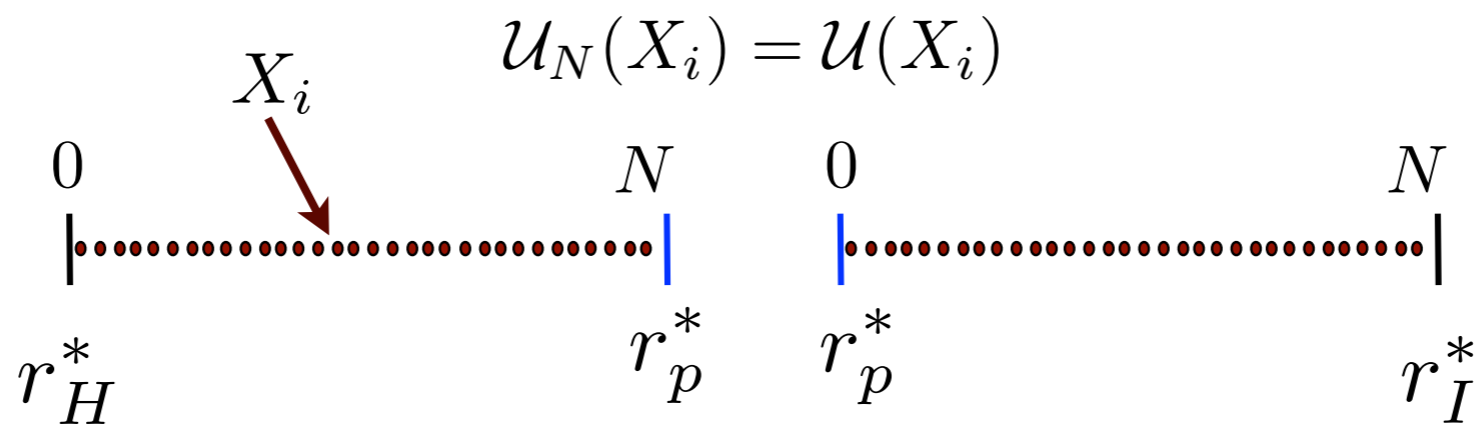
The Particle-without-Particle Scheme

- To implement the PwP scheme numerically we use **PSCM**. Each subdomain is discretised with a number N of collocation points of a Lobatto-Chebyshev grid:



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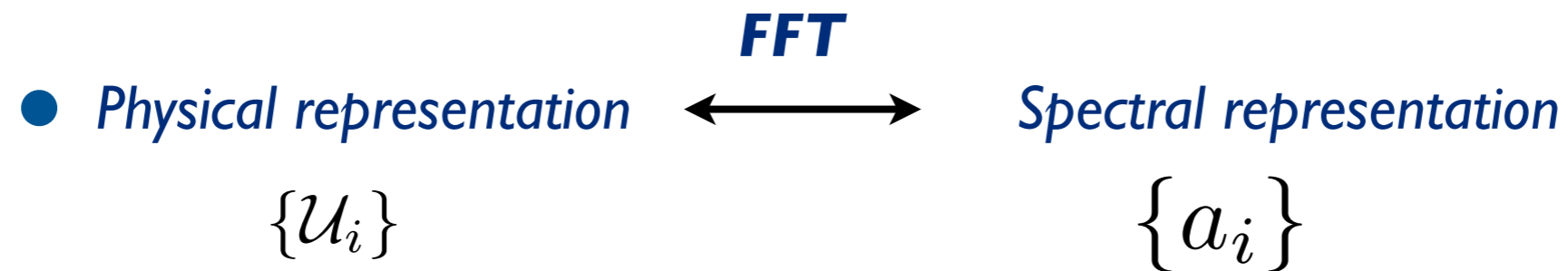


To prevent incoming signals from outside the physical domain

$$\begin{aligned} \phi^{\ell m}(t, r_H^*) - \varphi^{\ell m}(t, r_H^*) &= 0 = U^{\ell m}(t, r_H^*) & r_H^* &\rightarrow -\infty \\ \phi^{\ell m}(t, r_I^*) + \varphi^{\ell m}(t, r_I^*) &= 0 = V^{\ell m}(t, r_I^*) & r_I^* &\rightarrow \infty \end{aligned}$$

The Particle-without-Particle Scheme

Employing a Chebyshev basis there are some paybacks:



- Differentiation is cheaper in the spectral domain:

$$\mathcal{U}'_N = \sum_{j=0}^N D_{ij} \mathcal{U}_j(X)$$

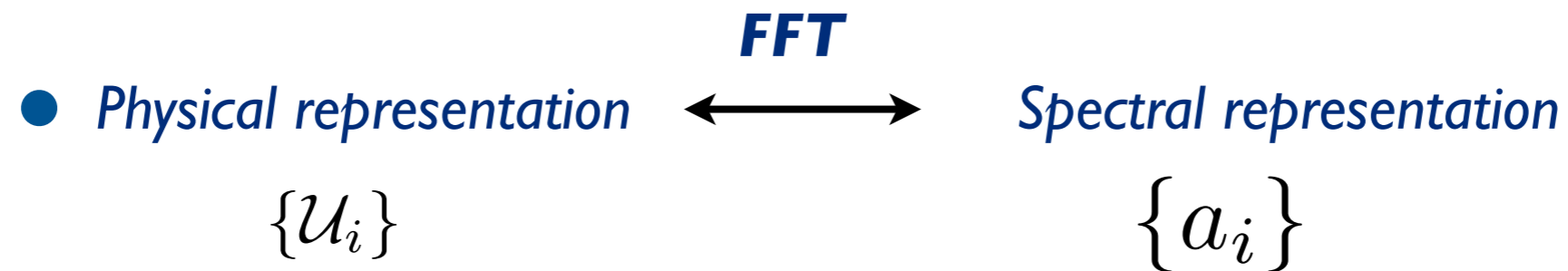
N^2 operations

$$\mathcal{U}'_N = \sum_{j=0}^N b_j T_j(X)$$

$\sim N \ln(N)$ operations

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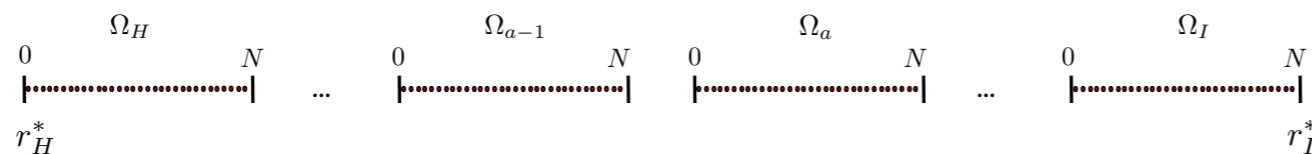
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$\sim \mathbf{N} \ln(\mathbf{N})$ operations

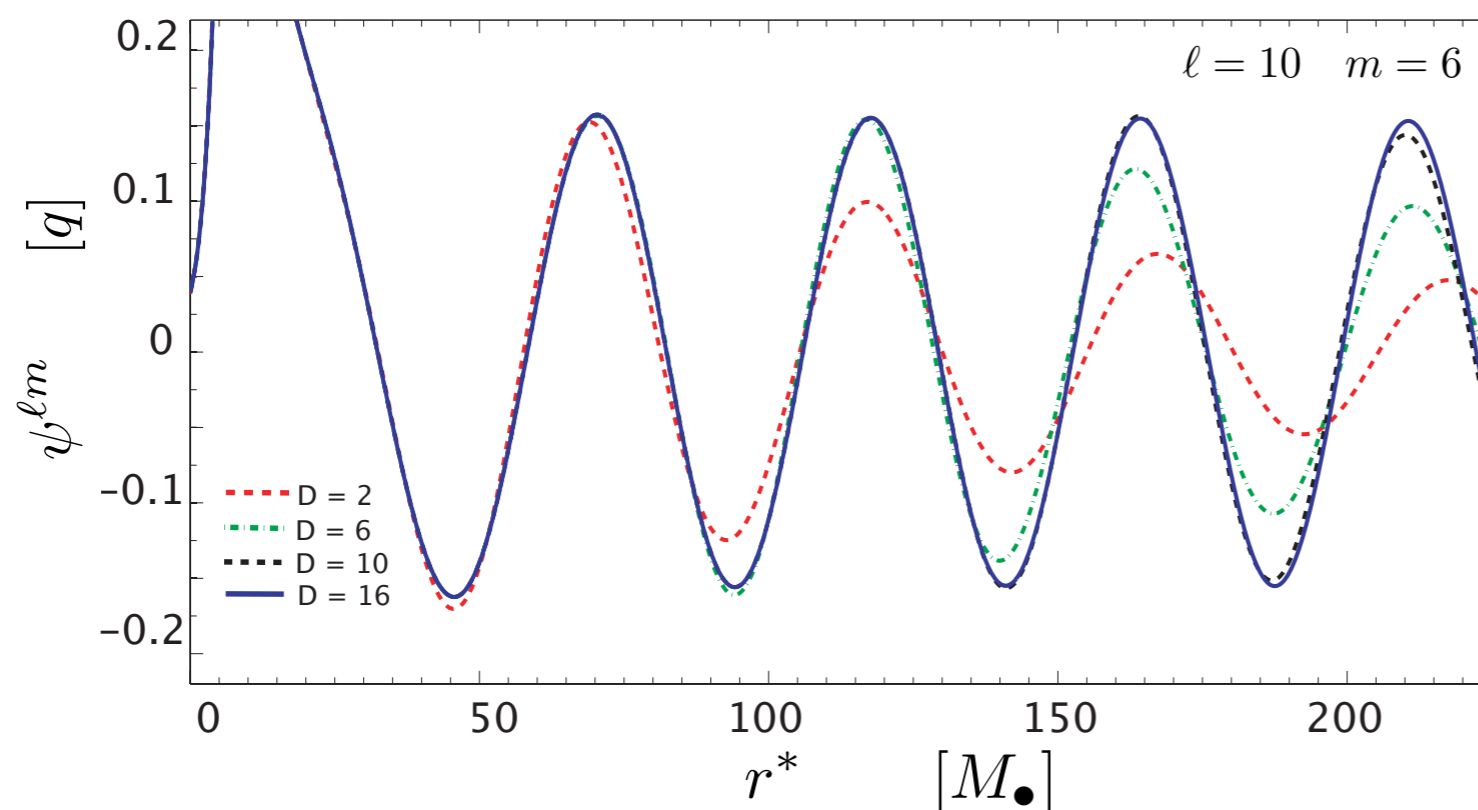
Our (smooth) solutions converge exponentially with \mathbf{N}

Multidomain flexibility

- Covering the spatial domain with a given number of subdomains (D) we improve the field resolution with a relatively small N .



$$\Delta r_a^* = r_N^* - r_0^*$$

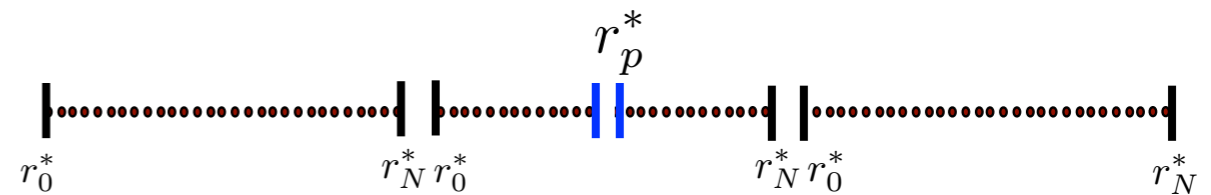
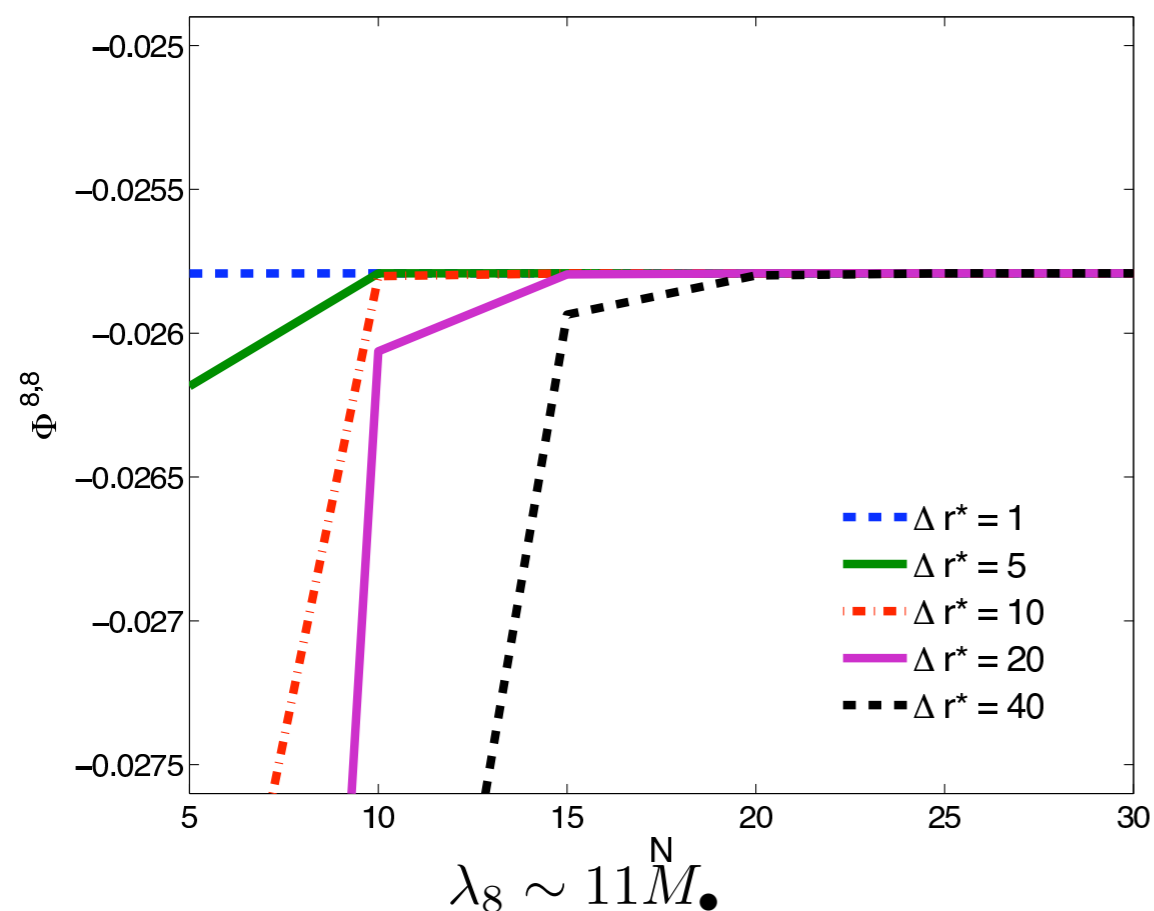


$$\Delta r^* = 50 M_\bullet \quad N = 50$$

$$\text{Resolution} = \left\{ \begin{array}{l} D \uparrow, \quad N \downarrow \\ D \downarrow, \quad N \uparrow \end{array} \right.$$

Advantages of the Multidomain Framework

- Different harmonic modes need different resolution
- We adjust the size of the subdomain around the particle location to the smaller mode wavelength



$$\Delta r^* = r_N^* - r_0^*$$

$$\Delta r^* \sim \lambda_m$$

$$\rho \sim \frac{\Delta r^*}{N}$$

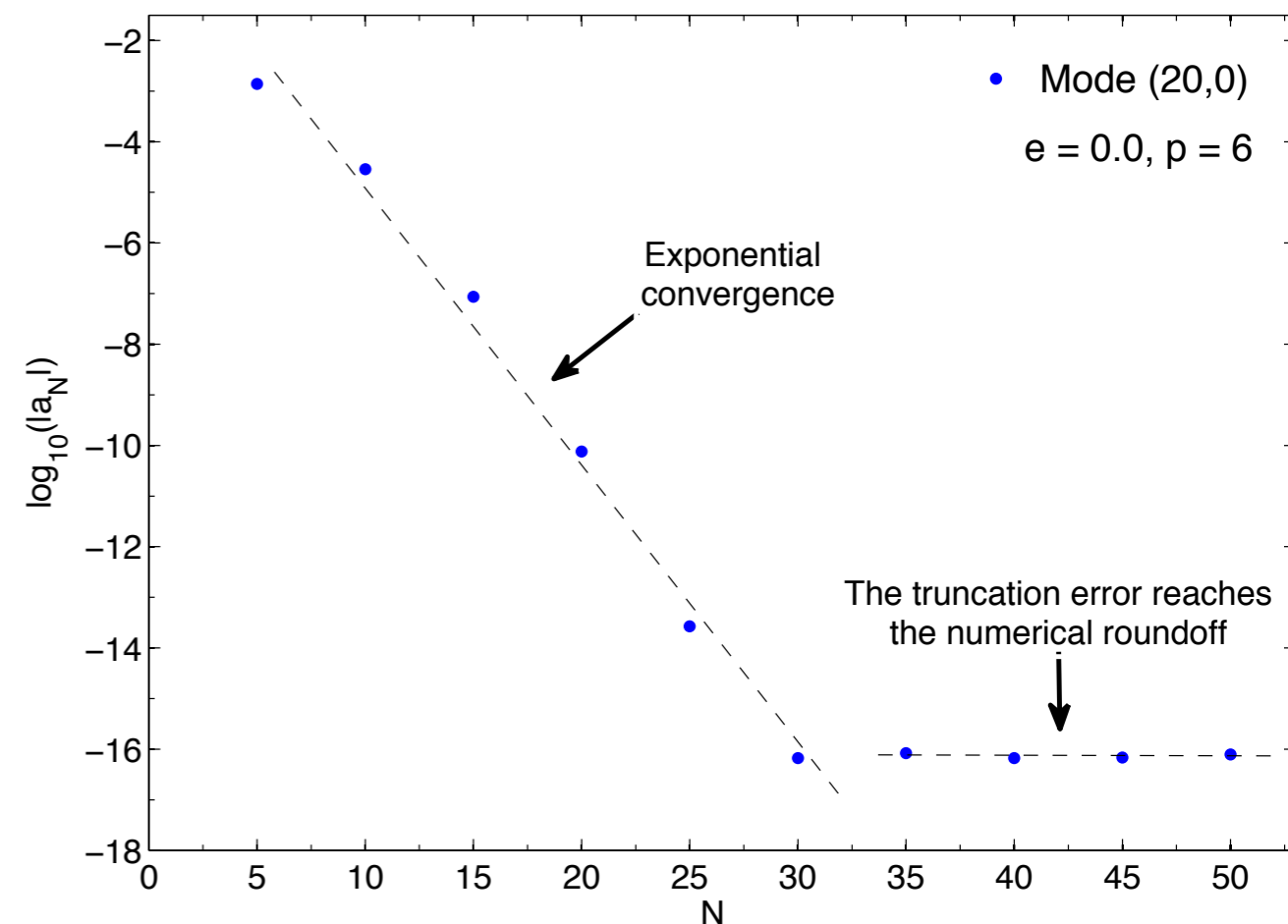
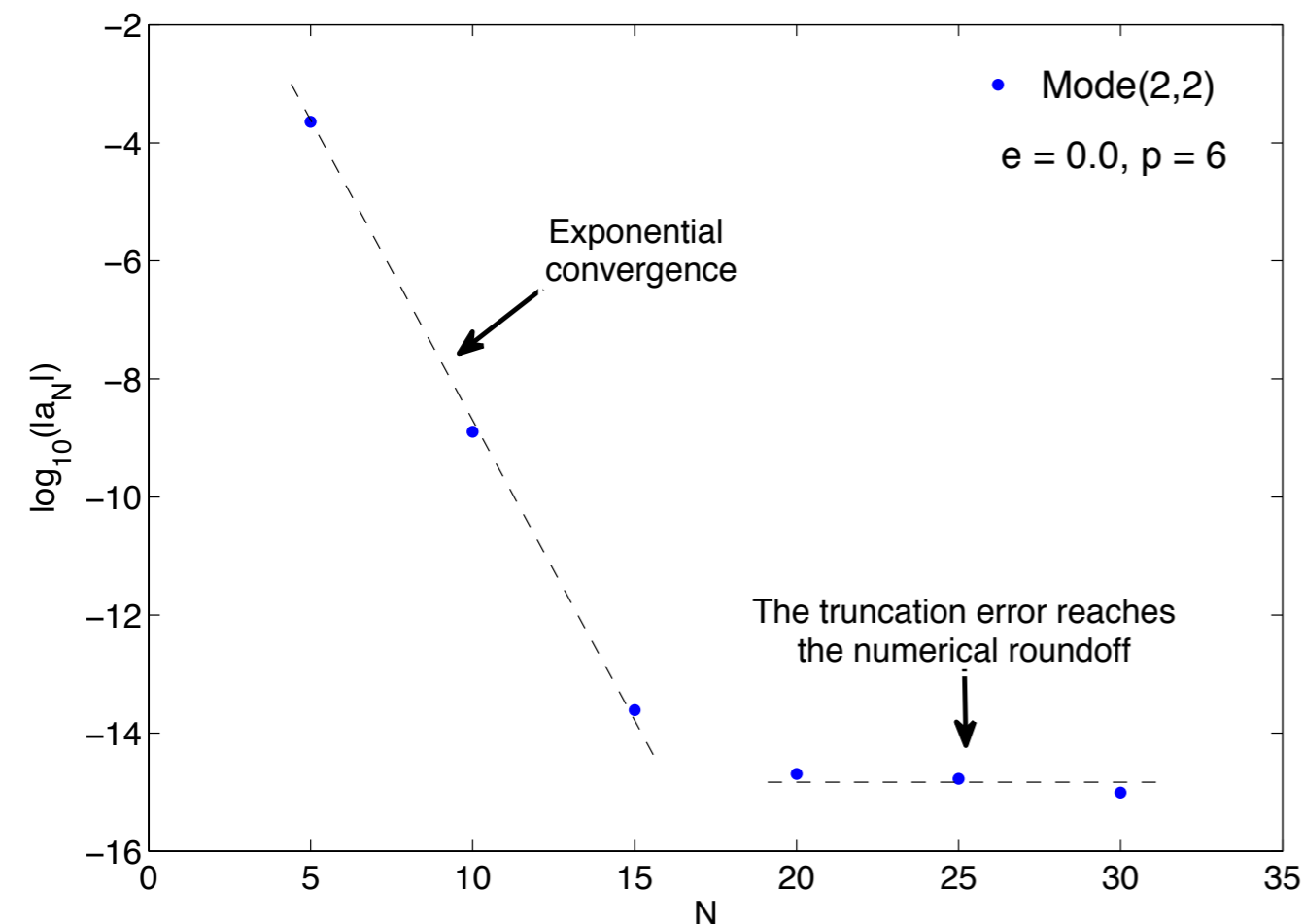
[Canizares & Sopuerta (2011)].

Convergence Test (circular case)

The dependence of the truncation error ($\sim |a_N|$) with respect increasing numbers of collocation points, N , give us an estimation of the exponential convergence of the code: e^{-N}

$$(\ell, m) = (2, 2)$$

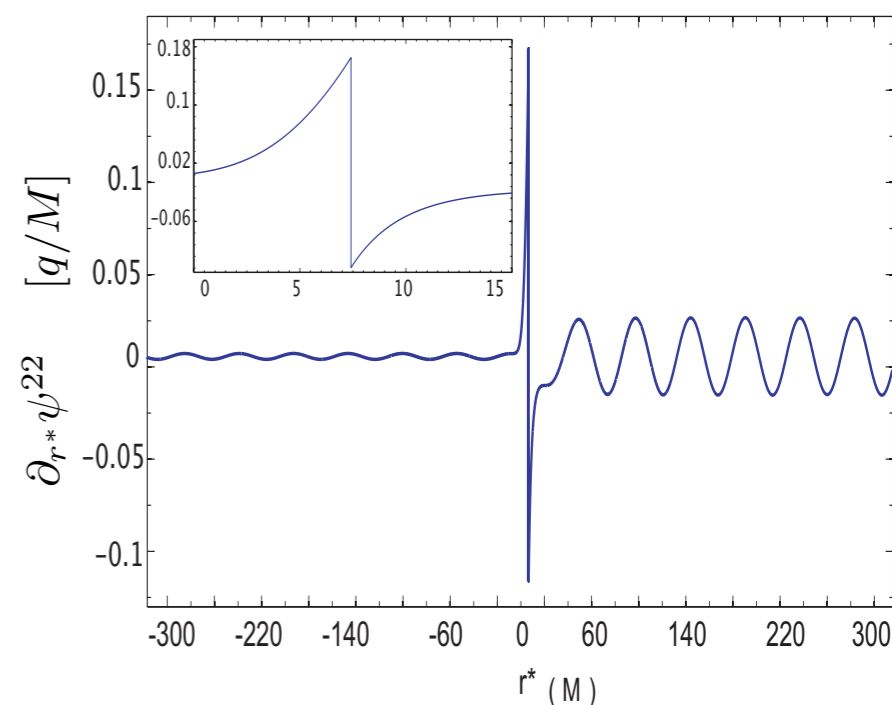
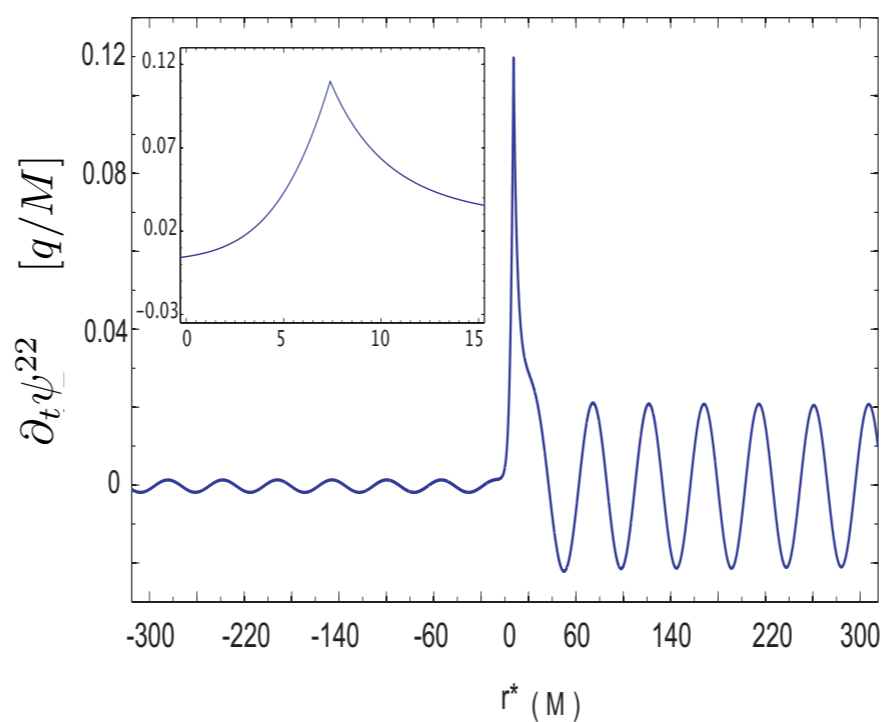
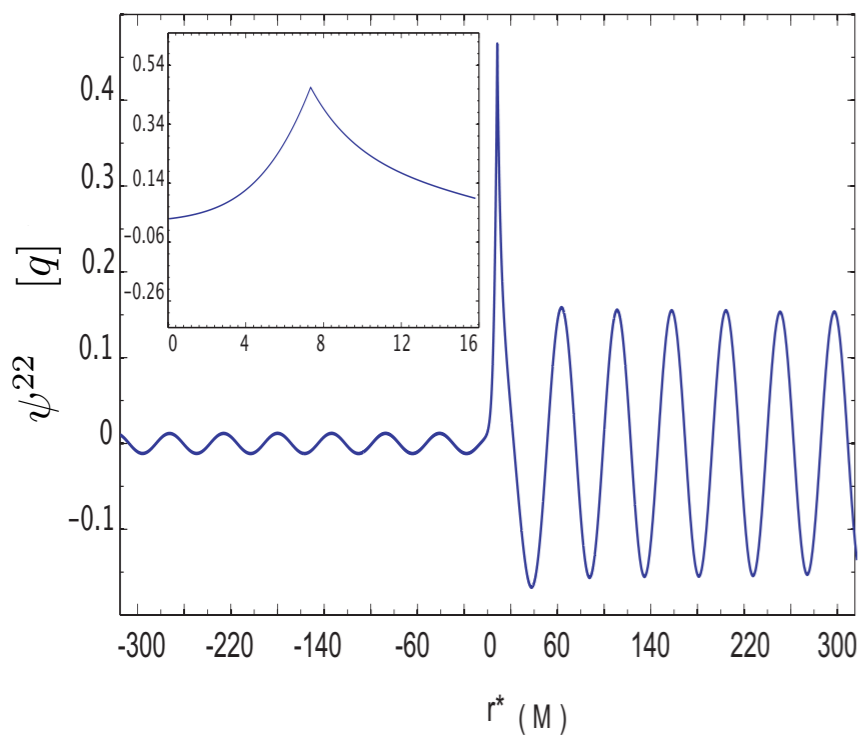
$$(\ell, m) = (20, 0)$$



[Canizares & Sopena (2009)] [Canizares & Sopena (2011)].

The Particle-without-Particle Scheme

Snapshots from the Circular case ($D=12, N=50$)



$$[\psi^{\ell m}]_p = 0,$$

$$[\partial_t \psi^{\ell m}]_p = 0$$

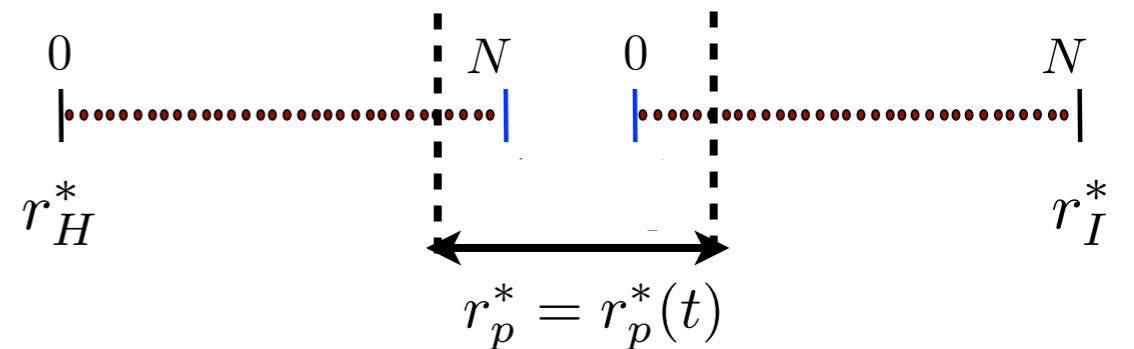
$$[\partial_{r^*} \psi^{\ell m}]_p = A^{\ell m}$$

[Canizares & Sopuerta (2009)]

The Particle-without-Particle Scheme

From circular to eccentric orbits:

- The key point of the PwP method is to keep the particle at the interface between subdomains:



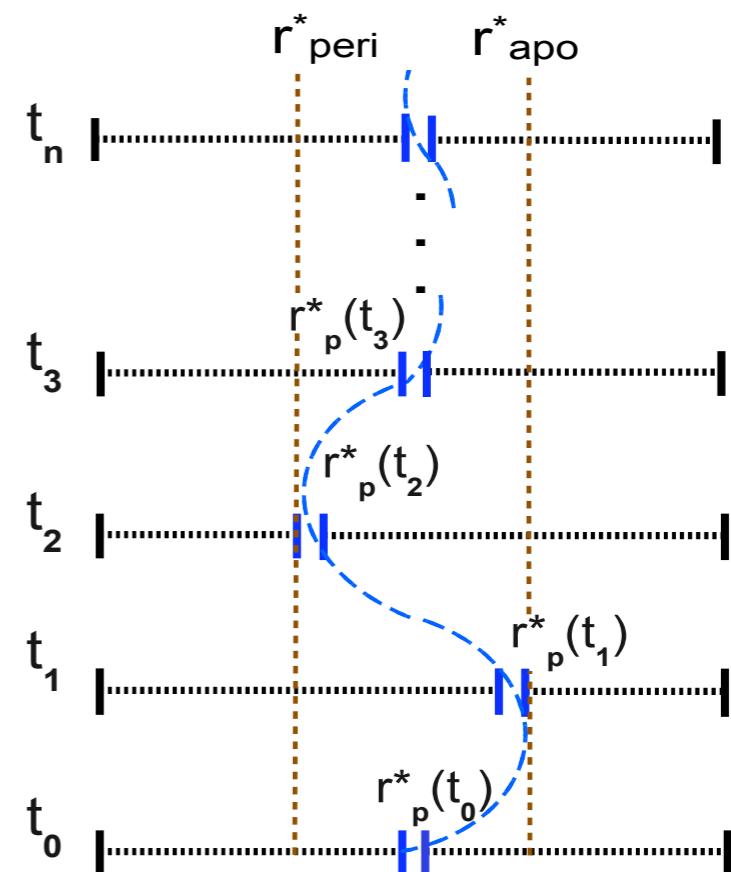
- For eccentric orbits we use a time dependent linear mapping between the physical and spectral domains.

$$r_p^* = r_p^*(t) \quad [\psi^{\ell m}]_p = 0 ,$$

$$[\partial_t \psi^{\ell m}]_p = - \frac{\dot{r}_p^* S^{\ell m}}{(1 - \dot{r}_p^{*2}) f(r_p)} ,$$

$$[\partial_{r^*} \psi^{\ell m}]_p = \frac{S^{\ell m}}{(1 - \dot{r}_p^{*2}) f(r_p)}$$

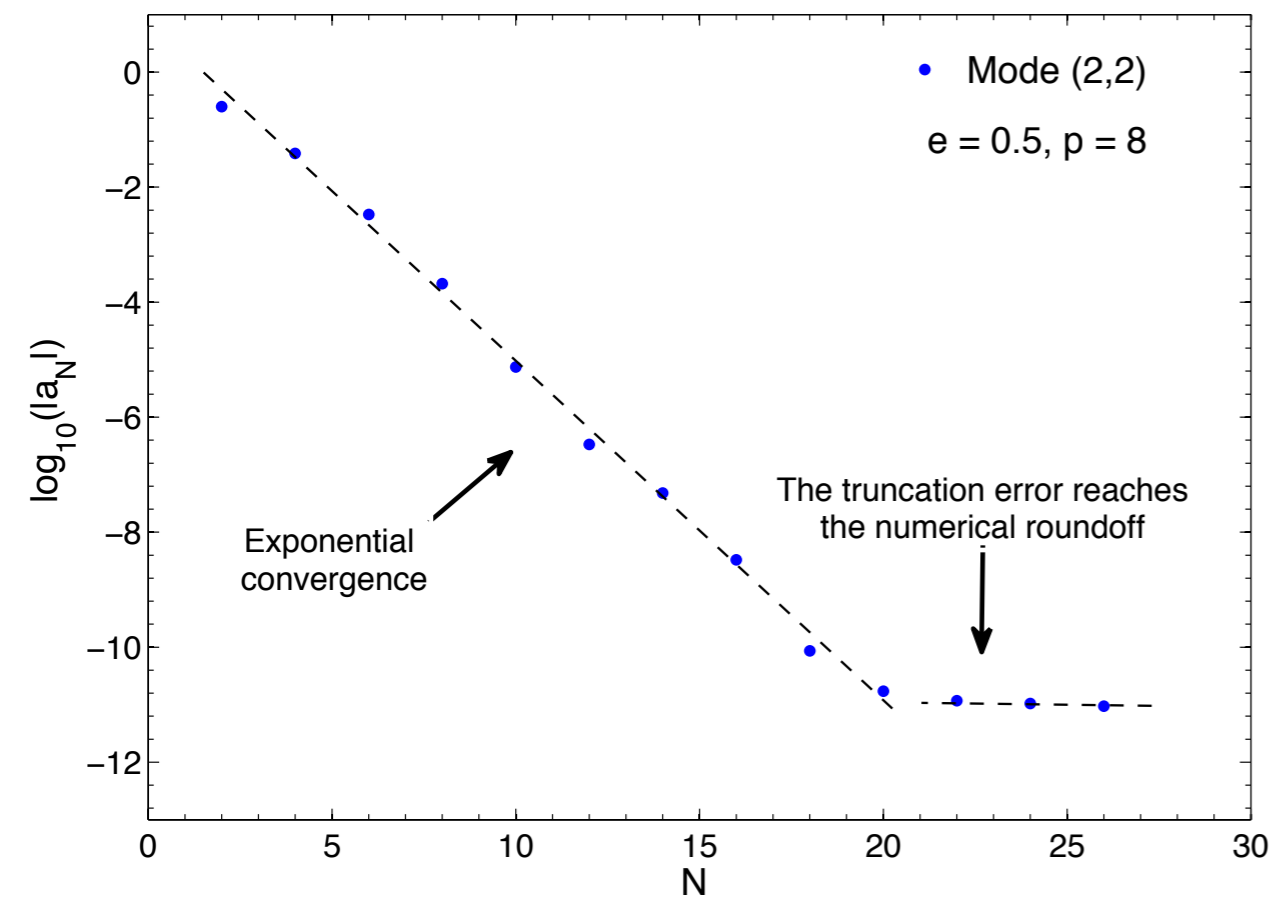
[Canizares, Sopena & Jaramillo (2010)].



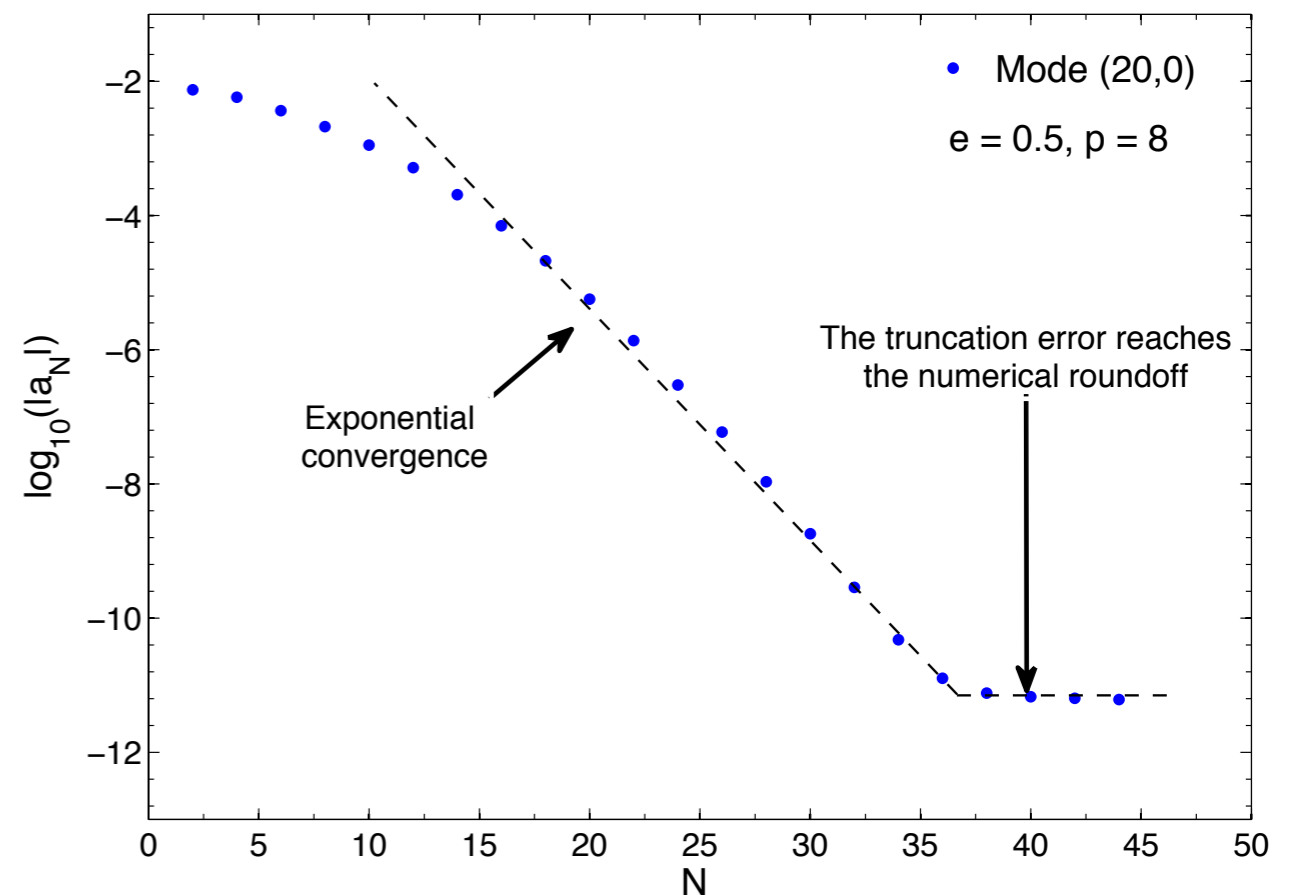
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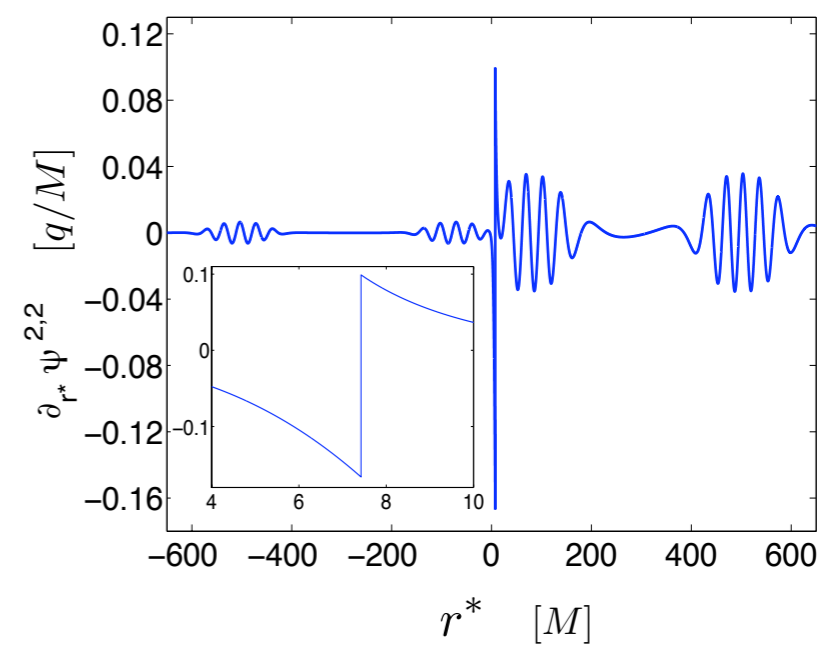
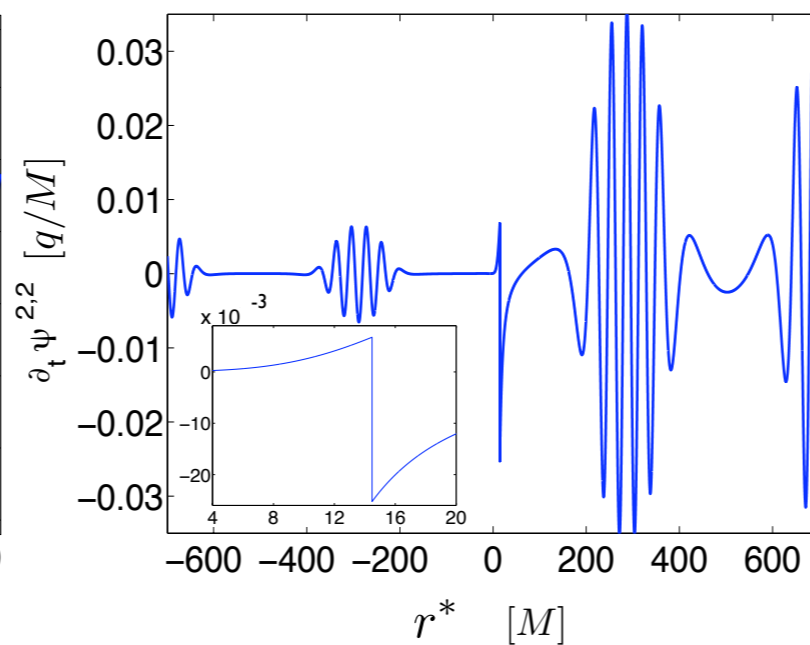
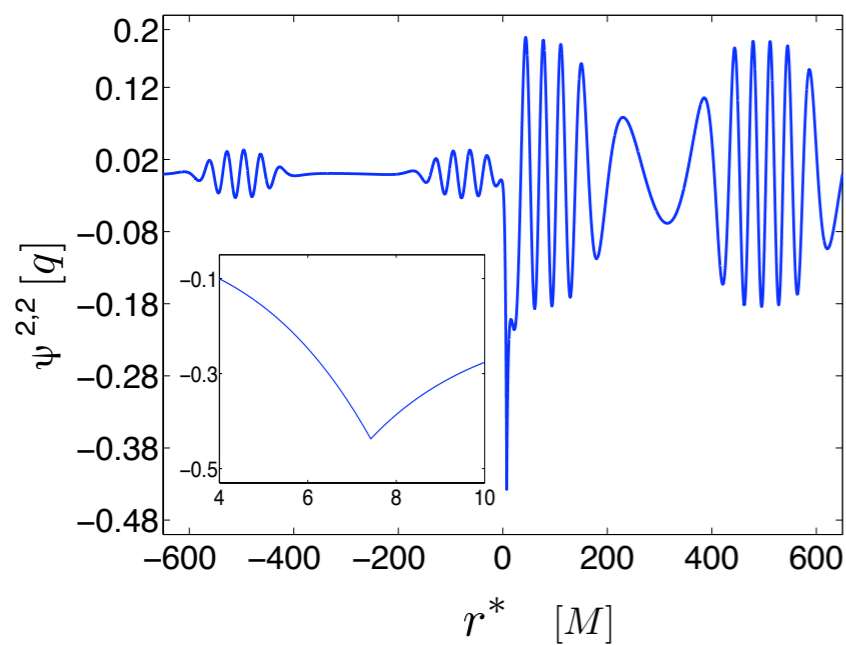
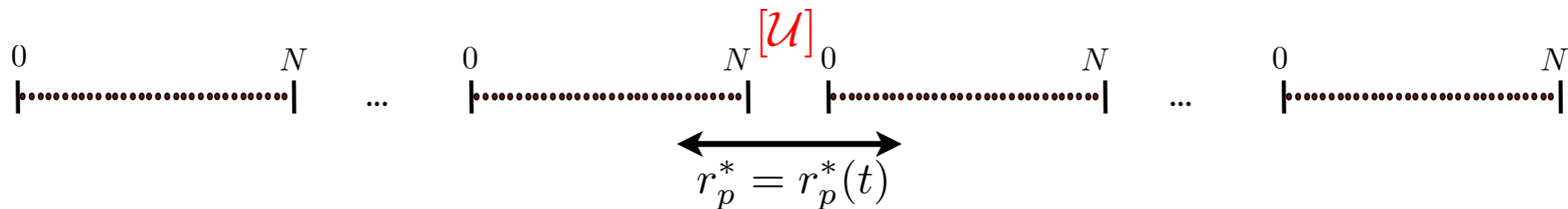
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[Canizares & Sopena (2011)].

The Particle-without-Particle Scheme

Snapshots from the Eccentric ($e=0.5, p=7.1$) case ($D=10, N=100$) $(\ell, m) = (2, 2)$



$$[\psi^{\ell m}]_p = 0, \quad [\partial_t \psi^{\ell m}]_p = -\frac{\dot{r}_p^* A^{\ell m}}{(1 - \dot{r}_p^{*2})}, \quad [\partial_{r^*} \psi^{\ell m}]_p = \frac{A^{\ell m}}{(1 - \dot{r}_p^{*2})}$$

[Canizares, Sopuerta & Jaramillo (2010)].

The Particle-without-Particle Scheme

Results for the self-force components:

- We have employed $\ell_{max} = 17$, $D = 10$, $N = 100$ and $\Delta r^* = 2 - 5M$.

The Self-force values have been obtained near the pericenter. [Canizares, Sopuerta & Jaramillo (2010)].

- We compare our results with posterior ones obtained in the frequency-domain

[Warburton & Barack (2010)]

| (e, p) | \mathcal{F}_α | PwP | Frequency-Domain | Relative Difference |
|------------|-----------------------------------|-----------------------------|----------------------------|-----------------------|
| (0.1, 6.3) | $\frac{M_\bullet^2}{q} \Phi_t^R$ | $4.517\ 196 \cdot 10^{-4}$ | $4.517\ 994 \cdot 10^{-4}$ | 0.01% |
| | $\frac{M_\bullet^2}{q} \Phi_r^R$ | $2.125\ 049 \cdot 10^{-4}$ | $2.125\ 7 \cdot 10^{-4}$ | 0.03% |
| | $\frac{M_\bullet}{q} \Phi_\phi^R$ | $-6.204\ 083 \cdot 10^{-3}$ | $-6.20\ 401 \cdot 10^{-3}$ | $3 \cdot 10^{-5}\%$ |
| (0.3, 6.7) | $\frac{M_\bullet^2}{q} \Phi_t^R$ | $7.698\ 048 \cdot 10^{-4}$ | $7.177\ 3 \cdot 10^{-4}$ | 0.25% |
| | $\frac{M_\bullet^2}{q} \Phi_r^R$ | $3.63\ 3926 \cdot 10^{-4}$ | $3.632\ 2 \cdot 10^{-4}$ | 0.04% |
| | $\frac{M_\bullet}{q} \Phi_\phi^R$ | $-9.040\ 222 \cdot 10^{-3}$ | $-9.0402\ 1 \cdot 10^{-3}$ | $1.5 \cdot 10^{-5}\%$ |
| (0.5, 7.1) | $\frac{M_\bullet^2}{q} \Phi_t^R$ | $1.233\ 071 \cdot 10^{-3}$ | $1.233\ 1 \cdot 10^{-3}$ | 0.015% |
| | $\frac{M_\bullet^2}{q} \Phi_r^R$ | $5.612\ 209 \cdot 10^{-4}$ | $5.617\ 9 \cdot 10^{-4}$ | 0.1% |
| | $\frac{M_\bullet}{q} \Phi_\phi^R$ | $-1.268\ 560 \cdot 10^{-2}$ | $-1.2685\ 7 \cdot 10^{-2}$ | $6.1 \cdot 10^{-4}\%$ |

Summary

- *The PwP scheme provides accurate and efficient self-force computations in $(1+1)$*
- *It is a robust method suitable to deal with generic EMRI orbits.*
- *We are working to extend the PwP scheme to $2+1$ computations*