## Increasing Fisher Information using Moving-Mesh Reconstruction

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Reconstruction techniques are commonly used in cosmology to reduce complicated nonlinear behaviour to a more tractable linearized system. We study the Moving-Mesh algorithm which is able to consistently compute the displacement field from non-linear density fields. To quantify the algorithm's ability to reconstruct linear modes, we study the Fisher information present in 130 N-body simulations before and after reconstruction. We find that the linear scale is pushed to  $k \simeq 0.3$  h/Mpc after reconstruction. We furthermore find that the non-linear plateau of the cumulative Fisher information is increased by a factor of  $\sim 50$  after reconstruction, from  $I \simeq 2.5 \times 10^{-5}/(\mathrm{Mpc/h})^3$  to  $I \simeq 1.3 \times 10^{-3}/(\mathrm{Mpc/h})^3$  at  $k \simeq 2.6$  h/Mpc. This result includes the decorrelation between initial and final fields, which has been neglected in some previous studies artificially improving their performance. We expect this technique to be beneficial to problems such as baryonic acoustic oscillations and cosmic neutrinos that rely on an accurate disentangling of nonlinear evolution from underlying linear effects.

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### I. INTRODUCTION

Two-point statistics provide a complete description of Gaussian density fields and can be computed efficiently even for large data sets. However, non-linear gravitational evolution leads to highly non-Gaussian matter distributions which require higher order statistics to fully characterize. Such statistics are computationally expensive and can be challenging to relate to cosmological parameters. To mitigate these difficulties, it is common to transform the matter field in a way that hopefully reduces non-Gaussianity. For example, Gaussianization transforms have been used to make the logarithmic distribution more Gaussian [1, 2] and Wavelet Non-Linear Wiener filters have been used to separate Gaussian and non-Gaussian components of the density field [3–5].

The success of techniques can be quantified by computing the Fisher information present in the power spectrum before and after reconstruction. Rimes and Hamilton [6] were the first to study the Fisher information in the non-linear matter power spectrum calculated from N-body simulations. They found that the information has a plateau on translinear scales ( $k \simeq 0.2-0.8 \text{ h/Mpc}$ ) due to strong coupling of Fourier modes. Qualitatively, this means that the power spectrum on these scales gives little additional information. However, Harnois-Déraps et al. [5] computed the Fisher information for various Gaussianization methods (and combinations of methods)

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and found that while mode coupling is reduced, there is not necessarily an improvement in the cross correlation between the initial Gaussian density field and the final non-linear one.

In studies of Baryon Acoustic Oscillations (BAO), density fields are subjected to reconstruction which partially inverts non-linear evolution by applying the negative Zel'Dovich displacement field [7]. The linear density field is typically estimated via Lagrangian perturbation theory (LPT) using the linear Zel'Dovich displacement  $-\nabla_q \cdot \Psi(q)$  with respect to initial coordiantes q [8]. Recently, Zhu et al. [9] described how to use the Moving-Mesh algorithm (MM), first described in [10, 11], to efficiently estimate  $\Psi(q)$  from non-linear density fields. They further showed that even though shell-crossing and vorticity are not recovered, linear density modes are still recovered up to scales relevent to the BAO.

In this paper, we compute the Fisher information recovered after using this reconstruction scheme on 130 independent N-body simulations. The paper is organized as follows. In §II, we briefly describe the computation of the displacement potential using MM algorithm. In §III, we describe the simulations, implementation of the reconstruction and compare the power spectra and cross correlations before and after reconstruction. In §IV, we futher compute the correlation matrix and Fisher information before and after reconstruction. Finally, in §V, we summarize our results and discuss the effectiveness of the reconstruction and its potential uses.

### II. RECONSTRUCTION ALGORITHM

In this section, we briefly review the MM algorithm; for a more complete description we refer the reader to [12]. The aim of the MM algorithm is to estimate the displacement of particles in Lagrangian coordinates from their final Eulerian position only. The general principle is to relate a particle's Eulerian coordinates,  $x^i$  to a curvilinear system,  $\xi^{\mu}$ , in which the number of particles per grid cell is approximately constant:

$$\rho\sqrt{g} = \text{const.},\tag{1}$$

where  $\sqrt{g} \equiv \det \left| e_{\mu}^{i} \right|$  is the volume element, and  $e_{\mu}^{i} = \partial x^{i}/\partial \xi^{\mu}$  is the coordinate transformation matrix. These coordinates are related via the so-called deformation, which we assume to be a pure gradient:

$$x^{i} = \xi^{\mu} \delta_{\mu}^{i} + \frac{\partial \phi}{\partial \xi^{\mu}} \delta^{i\mu}, \qquad (2)$$

and  $\phi$  is called the deformation potential. The physical picture is moving the girds based on  $\phi$  to adapt Eq. 1. Numerially, we can iteratively solve for the deformation potential via the diffusion equation,

$$\partial_{\mu}(\rho\sqrt{g}e_{i}^{\mu}\delta^{i\nu}\partial_{\nu}\dot{\phi}) = \Delta\rho, \tag{3}$$

where  $\Delta \rho = \bar{\rho} - \rho \sqrt{g}$  is the density difference after and before the grids are moved. A detailed discription of the analytical formulation can be found in [10, 11]. Eq. 3 can be solved through the use of the multigrid algorithm described in [10–12]. The displacement field is then given by

$$\Psi(\boldsymbol{\xi}) = \nabla \phi(\boldsymbol{\xi}),\tag{4}$$

and the reconstructed density field is given by

$$\delta_R(\boldsymbol{\xi}) = -\nabla \cdot \boldsymbol{\Psi}(\boldsymbol{\xi}) = -\nabla^2 \phi(\boldsymbol{\xi}). \tag{5}$$

# III. IMPLEMENTATION AND POWER SPECTRA

We use the CUBEP³M code [13] to run 140 simulations with a box size of 600 Mpc/h and 512³ particles. The initial conditions are computed using the transfer function given by CAMB [14] and then propagating the power back to z=100 with a linear growth factor. The Zel'dovich approximation is used to calculate the displacement and velocity fields of the particles. For these simulations, we use cosmological parameters  $\Omega_M=0.321,~\Omega_{\Lambda}=1.0-\Omega_m,~h=0.67,~\sigma_8=0.83,$  and  $n_s=0.96$ . Different random seeds are used to produce the initial conditions for different simulations so that they are independent of each other.

We use the Voronoi tessellation method to estimate the density contrast  $\delta_S = \delta \rho / \rho - 1$  from the particles, and then apply the MM reconstruction to these fields with a resolution of  $512^3$  cells. The reconstruction code solves the dispacement potentials iteratively until the root mean

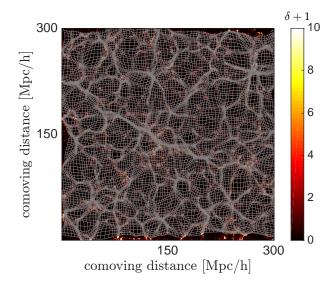


FIG. 1: The 2-D projection of one layer of the deformed grid of a sample N-body simulation is shown as curved white lines. The density  $\rho/\bar{\rho}=1+\delta$  is shown underneath.

square (rms) of the results drop from  $\sim 7.5$  to 0.20. For different simulation samples, a different number of iterations are required to get the results of the same rms. In total, 130 simulations converged to the target rms within 2000 iterations. A 2-D projection of one layer of the deformed grids and the original density field on the grids are given in Fig. 1. As expected, there is no grid crossing after reconstruction.

The cross power spectrum,  $P_{ab}(k)$ , is defined as

$$\langle \delta_a(\mathbf{k})\delta_b(\mathbf{k'})\rangle = (2\pi)^3 P_{ab}(k)\delta_{3D}(\mathbf{k} - \mathbf{k'}),$$
 (6)

where  $\delta_a$  and  $\delta_b$  are any density contrasts and  $\delta_{3D}$  is the three-dimensional Dirac delta function. We typically consider instead the dimensionless power spectrum,  $\Delta_{ab}^2(k)$ , defined as

$$\Delta_{ab}^2(k) \equiv \frac{k^3 P_{ab}(k)}{2\pi^2}.\tag{7}$$

In the left panel of Fig. 2, we show the matter auto power spectrum (a = b) of linear theory density fields  $(\delta_L)$ , from the simulation results  $(\delta_S)$  and after reconstruction  $(\delta_R = -\nabla^2 \phi)$ . For the simulation results, we use the average value of all 130 simulations and show  $1\sigma$  standard deviations as error bars.

To determine the correlation between fields, we compute the cross correlation coefficient  $r_{ab}(k) = P_{ab}/\sqrt{P_{aa}P_{bb}}$ . In the right panel of Fig. 2, we show  $r_{SL}$  and  $r_{RL}$ . We see that the reconstructed field is much more highly correlated with the linear field than the simulation field is. For comparison, we also plot the correlation coefficient of  $\delta_E$  and  $\delta_L$  from Yu et al. [15], where  $\delta_E(q) = -\nabla_q \cdot \Psi(q)$  is the negative divergence of the real non-linear displacement from simulaiton. Ideally, the MM algorithm aims to get the cross correlation  $r_{RL}$  close to  $r_{EL}$ . Even though  $r_{RL}$  decreases from  $r_{EL}$  in the non-linear regime, due to the fact that the MM reconstruction

cannot recover the cell-crossing and vorticity present on these scales, we find that linear modes are recovered successfully on scales  $k \simeq 0.05-0.3$  h/Mpc. Specifically, the scale at which r(k)=1/2 increases from  $k \simeq 0.2$  h/Mpc to 0.8 h/Mpc after reconstruction. In comparison with the results of Zhu et al. [12], we find the correlation coefficient falls off at slightly lower wavenumbers, which we attribute to using fewer particles per simulation.

### IV. FISHER INFORMATION CONTENT

Mathamatically, the Fisher information I of the initial scale invariant matter power spectrum, A, is defined as

$$I_A \equiv -\left\langle \frac{\partial^2 \ln \mathcal{L}}{\partial \ln A^2} \right\rangle,\tag{8}$$

in which  $\mathcal{L}$  denotes the likelihood [16]. In this paper, the word "information" and the symbol "I" both implicitly mean cumulative Fisher information of A. For Gaussian fluctuations, the likelihood depends on parameters only through the power spectrum P(k), so I can be written as

$$I = -\left\langle \sum_{k,k'} \frac{\partial \ln P(k)}{\partial \ln A} \frac{\partial^2 \ln \mathcal{L}}{\partial \ln P(k) \partial \ln P(k')} \frac{\partial \ln P(k')}{\partial \ln A} \right\rangle, \quad (9)$$

in which the angle bracket denotes the average of many realizations of the power spectrum [6].

Eq. 9 can be written in a simpler form in two aspects. First, we can simplify the derivative terms  $\partial \ln P(k)/\partial \ln A$ . For a given density field  $\delta_a$ , we can conveniently decompose it into linear and non-linear components

$$\delta_a(k) = b(k)\delta_L(k) + \delta_n(k), \tag{10}$$

in which  $\delta_L$  denotes the linear density field, b(k) is the bias and  $\delta_n(k)$  is defined such that the correlation  $\langle \delta_L(k)\delta_n(k)\rangle$  is zero. If we correlate  $\delta_a$  and  $\delta_L$ ,

$$\langle \delta_a(k)\delta_L(k)\rangle = b(k)\langle \delta_L(k)\delta_L(k)\rangle,$$
 (11)

we can solve for b as

$$b(k) = \frac{P_{aL}(k)}{P_{LL}(k)}. (12)$$

To find the non-linear term, we correlate  $\delta_a$  with itself,

$$\langle \delta_a(k)\delta_a(k)\rangle = b^2(k)\langle \delta_L(k)\delta_L(k)\rangle + \langle \delta_n(k)\delta_n(k)\rangle, \quad (13)$$

and find

$$P_{aa}(k) = b^{2}(k)P_{LL}(k) + P_{nn}(k).$$
 (14)

With the help of Eq. 12 and Eq. 14, we get

$$\frac{\partial \ln P(k)}{\partial \ln A} = \frac{P_{LL}(k)}{P_{aa}(k)} b^2(k) = r_{aL}^2(k). \tag{15}$$

The second step we can make is to simplify  $\partial^2 \ln \mathcal{L}/\partial \ln P(k)\partial \ln P(k')$  by utilizing the fact that its expectation value is the Fisher matrix. For linear fields, this is equal to the inverse of the covariance matrix which is diagonal with elements given by the number of modes in each bin (when considering k and -k as the same mode). We can extend this definition to non-linear fields, provided we take into account that the covariance matrix is no longer diagonal and invert it appropriately [6]. Thus, we can write the Fisher information in terms of matrix multiplication:

$$I(\langle k_n) = r^2(k)^{\mathrm{T}} \left[ C_{\text{norm}}^{-1}(k, k') \right]_{\langle k} r^2(k'),$$
 (16)

where  $C_{norm}$  is the normalized covariance matrix defined as

$$C_{\text{norm}}(k, k') = \frac{Cov(k, k')}{\langle P(k) \rangle \langle P(k') \rangle}, \tag{17}$$

r is the mean cross correlation of a given density field with linear one and the subscript  $< k_n$  indicates the matrix is set to zero for modes  $k, k' > k_n$ . The elements of the covariance matrix are defined as

$$\operatorname{Cov}\left(k,k'\right) \equiv \frac{\sum_{i,j=1}^{N} \left[P_{i}\left(k\right) - \left\langle P\left(k\right)\right\rangle\right] \left[P_{j}\left(k'\right) - \left\langle P\left(k'\right)\right\rangle\right]}{N-1},$$
(18)

where N is the total number of simulations and angle brackets are values averaged over all simulations.

The cross-correlation coefficient matrix, or for short the correlation matrix, is defined as:

$$\operatorname{Corr}(k, k') = \frac{\operatorname{Cov}(k, k')}{\sqrt{\operatorname{Cov}(k, k) \operatorname{Cov}(k', k')}}, \quad (19)$$

and represents the correlation between different k modes. The corelation matrices for non-linear and reconstructed power spectra are shown in the upper-left and lower-right sections of Fig. 3. By definition, the correlation matrix is symmetric with unit diagonal allowing us to overlay the two matrices. For the non-linear case, the correlation matrix is almost diagonal in the linear regime,  $k \lesssim 0.07 \text{ h/Mpc}$ . The off-diagonal elements are produced by strong mode coupling on non-linear scales and the super-survey tidal effect which is small on linear scales but dominates in the weakly non-linear regime [17]. The correlation matrix for the non-linear power spectra has a few negative elements (Corr  $\gtrsim -0.18$ ), which should vanish with more simulations [18]. For the reconstructed correlation matrix, the linear regime expands up to  $k \simeq 0.3$ h/Mpc. However, the number and magnitude of negative off-diagonal elements also increases (Corr  $\gtrsim -0.48$ ).

The cumulative Fisher information is proportional to the volume. We plot the cumulative Fisher information per unit volume of the simulation, linear and reconstructed power spectra in the left panel of Fig. 4. The Fisher information of the linear power spectra is equal to the number of k modes within the shell in wave space,  $N_k$ . We find that the Fisher information of the non-linear power spectra drops from the linear one at  $k \simeq 0.05$ 

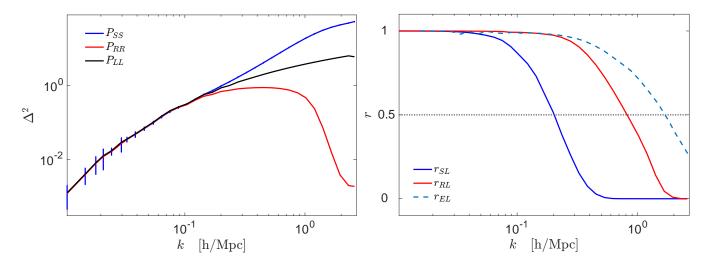


FIG. 2: Left. The dimensionless power spectrum computed via linear theory (black), the mean value of 130 N-body simulations with  $1\sigma$  error bars (blue), and reconstruction of the simulations (red). Right. The cross correlation function between simulation and linear densities  $r_{SL}$  (blue), MM reconstructed and linear densities  $r_{RL}$  (red), and E-mode reconstruction  $r_{EL}$  (dashed blue) from Yu et al. [15].

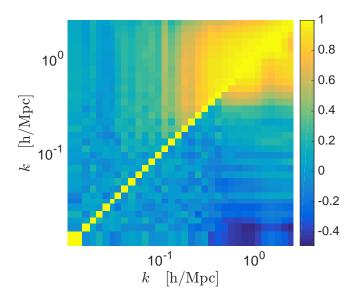


FIG. 3: The correlation matrix from 130 non-linear power spectra (the upper-left elements) and reconstructed power spectra (the lower-right off-diagonal elements).

h/Mpc, and has a flat plateau in the non-linear regime, with a saturated value of  $I \simeq 2.5 \times 10^{-5}/(\mathrm{Mpc/h})^3$ . This indicates that there is nearly no independent information in the non-linear regime. However, the information curve of the reconstructed power spectra keeps increasing roughly the same as the linear information until  $k \simeq 0.3$  h/Mpc, and reaches a value of  $I \simeq 1.3 \times 10^{-3}/(\mathrm{Mpc/h})^3$  at  $k \simeq 2.6$  h/Mpc, up by a factor of 50. This means that the MM reconstructed method can strongly recover the lost information within these scales. We compare the Fisher information given by the MM reconstruction method with the logarithmic density mapping method [2] as an example to illustrate its strength. We find that MM reconstruction gives over 10 times more information

than it. In some papers, the cross correlation  $r^2$  terms are set to unity in Eq. 16, which artificially increases the information. For comparison, we plot this case in the right panel of Fig. 4. We see that the logarithmic density mapping information is much higher, but only because it is not well correlated with the initial conditions.

Another way to quantify the non-linear scale is via the plateau's linear equivalent scale. That is, the wavenumber  $\kappa$  for which the linear information is the same as the non-linear plateau value, which we take to be at  $k \simeq 2.6$  h/Mpc. We find that the MM reconstruction increases  $\kappa$  from 0.15 to 0.4 h/Mpc whereas the logarithmic density mapping method only increases it to 0.19 h/Mpc.

### V. CONCLUSION

The MM reconstruction method successfully recovers the lost linear information on mildly non-linear scales and increases the saturated information from  $I \simeq 2.5 \times 10^{-5}/((\mathrm{Mpc/h})^3)$  to at least  $I \simeq 1.3 \times 10^{-3}/(\mathrm{Mpc/h})^3$ . The result is better than previous methods, e.g. [2–4, 19], and may improve further as the correlation coefficient between the reconstructed and linear fields will increase with higher resolution simulations [12]. This successful result on cold dark matter density fields provides strong motivation to adapt the MM reconstruction scheme to other cosmological fields such as biased tracers like halos and other matter components like baryons and neutrinos.

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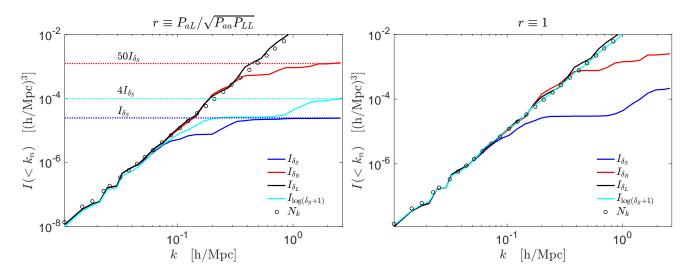


FIG. 4: Left. The cumulative Fisher information (solid lines) per unit volume as a function of wavenumber. The blue line corresponds to the non-linear density fields, the red line corresponds to the the reconstructed density fields, the dark line corresponds to the linear density fields, the cyan line corresponds to the logarithmic density mapping, and the circles are the cumulative number of k modes. Dotted horizonal lines indicate the value of the Fisher information at  $k \simeq 2.6$  h/Mpc. Right. Same as the left panel except with  $r \equiv 1$ . The black, blue and cyan lines match the results in [2, 6].

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