Caching and interpolated likelihoods: accelerating cosmological Monte Carlo Markov chains

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Abstract. We describe a novel approach to accelerating Monte Carlo Markov Chains. Our focus is cosmological parameter estimation, but the algorithm is applicable to any problem for which the likelihood surface is a smooth function of the free parameters and computationally expensive to evaluate. We generate a high-order interpolating polynomial for the log-likelihood using the first points gathered by the Markov chains as a training set. This polynomial then accurately computes the majority of the likelihoods needed in the latter parts of the chains. We implement a simple version of this algorithm as a patch (InterpMC) to CosmoMC and show that it accelerates parameter estimation by a factor of between two and four for well-converged chains. The current code is primarily intended as a “proof of concept”, and we argue that there is considerable room for further performance gains. Unlike other approaches to accelerating parameter fits, we make no use of precomputed training sets or special choices of variables, and InterpMC is almost entirely transparent to the user.

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Cosmological parameter values are typically estimated using Monte Carlo Markov Chains (MCMC). MCMC techniques are far more efficient than a brute force exploration of a parameter space with a realistic number of independent variables. Despite this, running Markov Chains for a broad range of parameter combinations — standard procedure when analyzing the cosmological implications of a major astrophysical dataset — remains computationally expensive. Consequently, algorithmic improvements that significantly increase the efficiency of this scheme without reducing its functionality are well worth exploring.

Schematically, MCMC parameter estimation begins with a model, or prior which has a set of free parameters $[1-5]$. A likelihood function (derived for a specific combination of datasets) returns the relative probability that the “observed sky” was produced by the prior with a specific set of parameter values. After picking an initial point in the parameter space, we compute the likelihood at a new set of parameter values. The chain will update to this new point with probability,

$$\alpha(\Theta_n, \Theta_{n+1}) = \min \left\{ 1, \frac{L(\Theta_{n+1})q(\Theta_{n+1}, \Theta_n)}{L(\Theta_n)q(\Theta_n, \Theta_{n+1})} \right\}$$

where $L(x)$ is the relative likelihood of parameter vector $x$ and $q(x,y)$ is the proposal density from $x$ to $y$. Within cosmology, CosmoMC is a canonical and widely used implementation of the algorithm [4].

A single likelihood can be computed in seconds, but a full set of chains requires the evaluation of many individual likelihoods. Computing $L$ is a nontrivial task since we must generate the CMB angular power spectrum (or $C_\ell$) corresponding to our chosen parameter vector. Further, improvements in both angular resolution and signal-to-noise in future datasets will require $C_\ell$ to be computed with increased precision and over a larger range of $\ell$ than currently necessary, sharply increasing the computational cost. Hence there is a strong need to accelerate the MCMC analysis of cosmological data.

A number of improvements to standard MCMC parameter estimation have been proposed to speed up or bypass the likelihood calculation. These include CMBFit [6], a polynomial fit to the WMAP 1-year likelihood, DASH [7], which uses a combination of precomputation and analytic approximation, and WARP [8], which combines interpolation with a careful choice of orthogonal variables [7] to accelerate the computation of the power spectrum. The
most mature packages are CosmoNet [9] which trains a neural network to provide likelihood values, and PICO [10, 11], which uses a large, precomputed training set.

MCMC estimation can be used with a vast range of problems. However, cosmological likelihoods — particularly those derived from the Cosmic Microwave Background (CMB) experiments — have two useful properties: they tend to be smooth functions\(^1\) of the input parameters \(x\) and evaluating them is computationally expensive. Consequently, we attempt to speed parameter estimates by caching computed values of the likelihood and then using interpolation to replace subsequent calls to the likelihood function within the MCMC code. The smoothness of the likelihood ensures that the interpolated likelihood will be a good approximation to the actual value. The computational overhead required by the caching and interpolation is small, compared to the cost of evaluating the likelihood directly, thus increasing the overall efficiency of the MCMC chains.

Crucially, we make no use of precomputation when constructing our interpolation. Our technique essentially works because the MCMC analysis itself “discards” enough information to reconstruct an interpolated likelihood; there is no need for a training set. This approach is straightforward computationally — we use a stock interpolation routine and make relatively small changes to CosmoMC itself. We implemented this algorithm as a patch to CosmoMC, dubbed InterpMC. Without any serious attempts to optimize the interpolation the runtime required for a typical parameter estimation is reduced by a factor of 2 to 4 for well-converged chains, with no degradation in the results. This improvement is nontrivial, although not as dramatic as that achieved by methods that rely on precomputation or analytic approximations. These can improve the runtime of MCMC code by an order of magnitude or more, but often at the cost of a great deal of extra work (whether analytic or computational), which must be performed beforehand. Further, InterpMC works for any combination of datasets and cosmological model, is almost entirely transparent to the user, and offers a good deal of scope for future improvement. The source is available as a patch file to CosmoMC.\(^2\)

The structure of this paper is as follows. In section 2 we describe our algorithm, and the details of its implementation. Performance metrics and tests used to validate the interpolation scheme are discussed in section 3. We test InterpMC with a variety of datasets (WMAP, ground based CMB, supernovae, BAO) and scenarios, including both the usual concordance cosmology, and models with curvature, neutrinos, running and tensors. Finally, in section 4 we summarize our work, and identify enhancements to InterpMC that could further improve its performance.

2 InterpMC: description and implementation

The key ingredient of InterpMC is a polynomial fit to computed likelihood values: this is a function of a given model’s free parameters. The interpolation data is gathered “on the fly”: the first 10–30% of the chains typically yield enough points to construct an accurate interpolation. We use likelihoods computed for both accepted and unaccepted points, and sets of chains run in parallel pool their interpolation data. Note that when using multiple datasets (e.g. BAO + WMAP7) we interpolate the full likelihood, rather than just the WMAP likelihood.

\(^1\)Exceptions to this rule certainly exist [12], but it will not be a particularly burdensome restriction.

\(^2\)http://easther.physics.yale.edu/interpmc.html - at this point the code is offered as “proof of concept”, rather than production-ready code, but it has proved to be robust in a wide variety of settings.
Figure 1. Flow chart illustrating how the interpolation is constructed.

We find that it is effective to fit the log-likelihood to an \( n \)-th order polynomial in the free parameters, rather than the likelihood itself. The likelihood is roughly Gaussian near the peak, and non-Gaussian corrections to the likelihood tend to be multiplicative, rather than additive. Consequently, while the likelihood and log-likelihood can both be expanded as polynomials, fitting to the log likelihood yields more stable results. Further, we exclude points from the interpolation which differ from the maximum log-likelihood value by more than a user-defined threshold, ensuring that the interpolation is not dominated by points that are infrequently visited by the Markov chains.

2.1 Constructing the interpolation

The \( n \)-th order interpolating polynomial for the log-likelihood contains all combinations of the \( P \) free parameters in the chains, up to \( n \)-th order. Consequently, the interpolation algorithm is required to estimate the parameters \( c_{ij...k} \) in the following polynomial

\[
\log L(x_1, \ldots, x_P) \approx c_0 + \sum_{i=1}^{P} c_i x_i + \sum_{i=1,j \geq i}^{P} c_{ij} x_i x_j + \sum_{i=1,k \geq j,j \geq i}^{P} c_{ijk} x_i x_j x_k + \cdots \tag{2.1}
\]

where the expansion will be truncated at order \( n \). The variables \( x_i \) are rescaled to have mean of zero and a standard deviation of unity, but are otherwise proportional to the “raw” variables in the chains.

As pointed out below, we make no claim that our fitting procedure is optimal, but do demonstrate that it is good enough to consistently improve the runtime of MCMC parameter estimates. The polynomial has \((P + n)!/P!n!\) unique coefficients, a number which obviously grows quickly with both \( n \) and \( P \). However, we were able to fit a 4-th order polynomial with
9 free parameters while substantially accelerating the corresponding parameter estimation. Within CosmoMC, the convergence of a set of chains is tested periodically. InterpMC uses these same checkpoints to determine whether the number of collected points (across all chains) passes a threshold value set by the user, which we typically chose to be three times the number of free parameters in the polynomial. When the threshold is exceeded, the chains share their databases of likelihoods. We first determine the peak likelihood encountered up to that point. The pooled points are then normalized and reduced to a subset whose log-likelihoods differ from the peak log-likelihood by less than some user-defined threshold. We then use an unweighted least-squares algorithm to fit both \( n-1 \)th and \( n \)th order polynomials to the cached likelihoods. By doing so, we can estimate the accuracy of the fit: if the two polynomials differ substantially, the interpolation is dominated by the highest order terms, and cannot be relied upon.\(^3\)

In addition to the free parameters in the chains, CosmoMC returns a number of derived parameters. Some of these, such as the age of the universe, only require the background FRW solution and can be trivially computed. Conversely, quantities such as the clustering parameter \( \sigma_8 \) or the alternative tensor:scalar ratio \( r_{10} \) require information obtained by running CAMB, which is precisely the step InterpMC is designed to avoid. Consequently, we also construct interpolations for these “slow” derived quantities. Given the quality of current data, these interpolations are extremely accurate since \( \sigma_8 \) and \( r_{10} \) vary more slowly than the likelihood, as functions of the free parameters. A flowchart outlining the use of the interpolation is shown in figure 1.

### 2.2 Using the interpolation

Once the polynomial coefficients have been calculated, the chains continue to run. As noted above, several criteria must be met for the interpolated likelihood to be used. First, the

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\(^3\)We use a version of Applied Statistics algorithm AS 174 for the fitting [13].
Figure 3. We plot the ratio of wallclock time to the number of accepted steps for ten different chainsets, for three different estimations. For a given model, the total number of steps varies by a factor of a few for a given model but the ratio of runtime to the number of steps is roughly constant. The models shown here are a) WMAP + ground based CMB for ΛCDM + w, b) WMAP+BAO with ΛCDM + r and c) WMAP + BAO + Supernovae with ΛCDM + n_{run}.

| Shorthand Variables (+7 Params) |
|---------------------------------|
| Reference | Concordance Only |
| Dark energy equation of state | w |
| Curvature | Ω_k |
| Tensor:scalar ratio | r |
| Running spectral index | n_{run} |
| Curvature + tensors | Ω_k + r |
| Tensors + running | r + n_{run} |
| Neutrino fraction | f_ν |

Table 1. The variables used for the trial runs.

interpolated likelihood must lie above the user-defined threshold, setting the maximum difference between the interpolated value and the peak value. Secondly, the n − 1th and nth order values are compared, and must agree to within a specified tolerance. When all these conditions are met, the chain obtains the log-likelihood from the nth order polynomial, and then accepts or rejects the corresponding step as usual. If these conditions are not satisfied, the full likelihood code is called. If the regular likelihood code is called, the computed point is added to the set of cached values, and the interpolating polynomials are re-estimated each time CosmoMC tests for convergence. Figure 2 summarizes the workings of INTERPMC.
3 Testing and results

To test InterpMC we ran chains for a number of parameter sets. As well as the standard concordance variables, we examined scenarios with a non-trivial neutrino sector, spatial curvature, tensors, and a running index, as listed in table 1. For each combination of parameters, we ran chains using only WMAP data, and WMAP in conjunction with other data, including supernovae, large scale structure [LSS] and Baryon Acoustic Oscillation [BAO], as listed in table 2. For each choice, we ran both interpolated and conventional chains, and compared the resulting runtimes and estimated parameter values. We have seven combinations of datasets, and eight combinations of variables, giving a total of 56 distinct estimations. In each case we found excellent agreement between the interpolated and non-interpolated chains.

MCMC estimates are inherently stochastic, so there is some variation in the runtimes for otherwise identical estimations, since the path taken by each chain through the parameter space is a random walk. However, we found that for a given combination of dataset and model, the ratio between the number of steps and the runtime is roughly constant, while the runtime itself varies by a factor of 2 or more. However, the ratio of runtime/attempted steps...
Figure 5. Histogram showing fraction of points in an MCMC chain set for which the likelihood can be obtained via interpolation, once enough data has been accumulated to construct the polynomial. Points whose interpolated log-likelihood differs by more than 8 (set by the cut parameter) are automatically computed from the full likelihood, after running CAMB. Points which survive this cut can almost all be obtained directly from the interpolation. This plot shows results for $\Lambda$CDM chains, using only the WMAP7 likelihood.

is an accurate measure of performance, as illustrated in figure 3, and we use this metric to evaluate performance gains associated with InterpMC. This constant is a function of the chosen dataset, model, and priors.

We worked with the January 2010 release of CosmoMC. We used the WMAP7 likelihood code; likelihoods for other datasets are supplied with CosmoMC. We started with the default settings for the initial proposal matrix with periodic updates enabled, and set MPI_Converge_Stop to 0.01. This is more conservative than the default value of 0.03 but yields smooth two dimensional parameter contours. This choice extends the runtime and thus increases the acceleration induced by the interpolation, since a greater fraction of the chains can use the interpolated likelihood. The interpolation typically commences well before the chains are close to convergence.

As well as comparing the parameter estimates with and without interpolation, we test the interpolation scheme by calling the regular likelihood and comparing the result to the interpolated value. The absolute error in the log-likelihood for the interpolated points was on the order of 0.01 – 0.05, with increasing error for values further removed from the peak likelihood. In figure 4 we plot the error versus the likelihood. Given that our parameter estimates overlap well with those returned by the stock version of CosmoMC, it is empirically clear that this error is not having a dramatic effect on the parameter estimates.

More quantitatively, looking at figures 4 and 5 we see that essentially all the accepted points in the chains have $\log L$ that differs by less than 10 from the maximum likelihood. Consequently, we do not need a particularly accurate estimate of $\log L$ for points far from the peak, since these points are almost always rejected. Conversely a “typical” interpolated point lies has a value of $\log L$ that differs by perhaps 3 or 4 from the peak value, and can be interpolated with an error, $\Delta \log(L) = \log L - \log L|_{\text{interp.}}$, of perhaps 0.02. Consequently,
the resulting error in $\mathcal{L}$ is

$$\mathcal{L} = e^{\Delta} \exp \left[ \log \mathcal{L}_{\text{interp}} \right]. \quad (3.1)$$

For the most majority of interpolated points, $|\Delta| < 0.025$, so the “exact” and interpolated likelihoods differ by a few percent. However, given that MCMC processes are intrinsically stochastic the interpolation error is small enough to be effectively unresolved within the parameter estimates below.

The user-supplied parameters that control the interpolation are added to the CosmoMC params.ini file, and are summarized in table 3. We set the $cut$ parameter to 8, so that interpolation is only used for points where the log-likelihood is within 8 of the maximum value. This encompasses most of the “peak” since it excludes points who relative likelihood is $e^8 \approx 3000$ times less than peak value. Secondly, we require that the likelihoods computed from the $n-1^{\text{th}}$ and $n^{\text{th}}$ order interpolations differ by no more than 20%. This is not a particularly stringent cut and rejects only a few points, although some of them very close to the peak. Finally, we used $n = 4$, so were comparing 3rd and 4th order interpolations. We show the fraction of calls to the likelihood function that are computed via interpolation in figure 5.

We ran chains for 56 different combinations of datasets and variables, using both InterpMC and the original CosmoMC code. We assessed the required runtime by computing the ratio of runtime / accepted point for each set of chains. Each estimation was run using 8 chains on an dual Intel “Xeon” cpu node (E5440@2.83 GHz). InterpMC does not add significantly to the MPI overhead, so CosmoMC continues to be “embarrassingly parallelizable”, and the interpolation routines take negligible amounts of runtime.

We ran with the parameters defined in the “stock” params.ini distributed with CosmoMC, apart from changing the convergence parameter to 0.01 so as to ensure that the 2 dimensional parameter constraints were smooth, as noted above. This allowed the chains to update the proposal matrix as they ran, which shortens the runtime unless the initially supplied covariance matrix is already close to optimal.

The performance of InterpMC is shown in table 4, and we see that the typical improvement in runtime is between 2 and 4. We look at models with up to 9 free parameters, and do not see an obvious correlation with between the performance gain and the total number of parameters. Figure 6 shows the performance of InterpMC as a function of the chosen set of free parameters, while while figure 7 shows the performance as a function of the datasets being employed for the estimation. Finally, looking at representative parameter estimates shown in figures 8 and 9 we see that the likelihood contours produced by CosmoMC and InterpMC...
| VARIABLE/DATASET | 7 Params | $\Omega_k$ | $r$ | $n_{\text{run}}$ | $\Omega_k + r$ | $r + n_{\text{run}}$ | $w$ | $f_{\nu}$ |
|-----------------|---------|---------|------|----------------|----------------|----------------|------|---------|
| WMAP7           | 3.5378  | 2.3794  | 3.5797| 2.8896        | 2.4169        | 2.4409        | 3.0976| 3.3019  |
| gCMB            | 3.3525  | 2.4504  | 3.5060| 3.0079        | 2.7707        | 2.1830        | 2.9112| 3.8758  |
| LSS             | 3.3577  | 2.7639  | 2.8640| 2.0883        | 2.5246        | 2.4409        | 1.8909| 3.5843  |
| LSS+gCMB        | 3.4446  | 2.9328  | 2.7487| 2.6698        | 2.7068        | 2.5653        | 2.7886| 3.6044  |
| LSS+gCMB+SN     | 3.3576  | 2.2650  | 2.3680| 2.2415        | 2.4163        | 2.8079        | 1.4848| 3.2155  |
| WMAP+BAO        | 3.5994  | 4.3205  | 3.1864| 2.4008        | 2.0084        | 2.0426        | 3.0084| 3.5202  |
| WMAP+BAO+SN     | 3.0137  | 3.1614  | 3.1986| 2.2648        | 2.0851        | 2.3197        | 3.4321| 3.3011  |

Table 4. Performance improvement for InterpMC, relative to standard CosmoMC for different choices of parameters and dataset.

InterpMC are effectively identical; the only difference between them is in the runtime required to produce them.

4 Discussion

We describe an approach to reducing the runtime required for cosmological parameter estimation, based on interpolating cached likelihood values computed with a single set of MCMC chains. An implementation of this scheme, InterpMC, is available as a patch to the standard MCMC package CosmoMC. This analysis focuses on cosmological parameter estimation, but the fundamental approach is applicable to other problems where the likelihood is a smooth function of the free parameters, and computationally intensive to evaluate.

We should ask why this approach works. Firstly, and unlike methods based on a precomputed training set (unless one has a very good idea of where to look, which is really only possible after an estimation has been performed), Markov Chains primarily evaluate points in regions of high likelihood, so our dataset is naturally weighted towards those points that will be frequently sampled during the estimation. Secondly, while MCMC estimates are much more efficient that a brute force search of the parameter space, they have no “memory” — their behavior is determined solely by the current location of the chain in parameter space. Any interpolation scheme, whether generated on the fly or via a precomputed training set is, to some extent, exploiting the global properties of the likelihood function. In particular, cosmological likelihoods are typically smooth functions of the free parameters we are seeking to estimate. This greatly facilitates the construction of an interpolating polynomial, but this information is not exploited by the MCMC algorithm itself. As a side note, we would point that cosmological parameter estimation already makes substantial use of interpolation: CAMB directly computes only a subset of the underlying Fourier modes and the $C_\ell$ values and then fits the intervening points. Turning off this interpolation significantly increases the runtime of CAMB. Consequently, InterpMC is extending the use of interpolation in cosmological parameter estimation, rather than introducing it for the first time.

As presently implemented, the efficiency gain yielded by InterpMC is typically between 2 and 4. Our primary goal here is simply to demonstrate the feasibility of this approach. We have deliberately employed conservative choices of the adjustable parameters in order to demonstrate that this approach works without careful tuning. Our experiments have shown that choosing more aggressive settings for the free parameters in InterpMC can noticeably increases the efficiency gain delivered by InterpMC without any modification to the underlying algorithm. Further, the interpolation requires a very small fraction of the total
Figure 6. Runtime per (attempted) step for InterpMC (triangle) and stock CosmoMC (square) runs versus the total number of steps, for each parameter set. Black lines connect runs performed with InterpMC and CosmoMC. The required number of steps varies stochastically even for otherwise identical runs, and there is no systematic difference between the number of steps required by InterpMC, relative to the stock code.

runtime, so even in the worst case scenario it cannot significantly slow CosmoMC, relative to a standard run.

We can see several specific algorithmic changes that may significantly improve the performance of InterpMC. Firstly, we currently fit to all combinations of all free parameters, up to order \( n \). However, many of these parameters are largely uncorrelated, and a more intelligent (but still automated) fitting procedure would focus on the subset of parameter combinations which make a nontrivial contribution to the fit, allowing the chains to begin
making use of interpolated likelihoods at an earlier point in the run. Separately, the current interpolation is expressed in terms of the (normalized) free parameters in the chains. However, we can always choose a basis in which the free parameters are uncorrelated at second order, so that the coefficient of $x_i x_j$ in the interpolating polynomial is proportional to $\delta_{ij}$. In this case, the higher order coefficients (e.g. $c_{ijk} x_i x_j x_k$) would reflect the couplings between variables whose covariance vanishes at second order, and will likely have a smaller number of nontrivial terms than the corresponding polynomial written in terms of the unrotated variables. Further, we could move to a scenario in which some combinations of variables are

Figure 7. Runtime per (attempted) step for InterpMC (triangle) and stock CosmoMC (square) runs versus the total number of steps, for each combination of datasets. Black lines connect runs preformed with InterpMC and CosmoMC.
Figure 8. We show 1D and 2D marginalized likelihood plots for the reference 7 parameter run using only WMAP, comparing results from an unmodified CosmoMC run, and results from InterpMC.

interpolated at higher order than others. Finally, while we have focused on the ability to avoid precomputation, it would be simple to save the computed likelihoods to “seed” a future run\(^4\), or to accelerate the subsequent computation of Bayesian evidence.

Our goal here has been to interpolate the likelihood in a way that is essentially invisible to the user and does not modify the underlying MCMC algorithm in any way. However,

\(^4\)This might be useful in situations where the likelihood computation was unchanged, but the allowed parameter ranges were altered, or the chains were to be extended for better convergence.
we have effectively demonstrated that we can construct an accurate interpolation to the likelihood for a wide range of cosmological datasets and models, while expending less computational effort than that which is required for a full parameter estimation. The interpolating polynomial effectively yields the functional form of the likelihood in a substantial volume surrounding the peak. Consequently, the existence of a closed-form algebraic expression for the likelihood may allow us to pursue parameter estimation techniques that make direct use of this information, beyond standard MCMC techniques.
It is clear that the computational cost of cosmological parameter estimation will continue to rise, particularly once data from Planck becomes available. Not only will this dataset include information at smaller angular scales than WMAP, the signal to noise will also be substantially improved, requiring a more accurate evaluation of the theoretical $C_\ell$, via CAMB. The computational cost of CAMB rises rapidly with the precision of the $C_\ell$ [11], so this problem will not be solved simply by relying on the continually improving price to performance ratio of computing hardware. Schemes relying on precomputation can be very fast at runtime, and even if the the production of the training set is computationally expensive this can be shared widely within the community. However, any training set may need to be regenerated when CAMB (or an equivalent code) is updated. Further, if one wishes to analyze a nonstandard set of independent variables there is no guarantee that the corresponding models can be expressed accurately in terms of the free parameters of the precomputed training set. These situations might include direct constraints on inflationary parameters (e.g. [14]), non-trivial power spectra, a complicated neutrino sector, or a nontrivial reionization history. On the other hand, if one does not need accurate parameter estimates (which require running long after the chains have completed burn-in) precomputation has significant advantages, especially if the likelihood can be evaluated with reasonable accuracy far from the peak.

To summarize, we have presented an initial implementation of an interpolation driven approach to accelerating MCMC cosmological parameter estimations, and shown that it can produce accurate results while reducing the resulting computational expenditure by at least a factor of two with a simple “proof of concept” implementation of this algorithm. Further, we have identified specific improvements to this approach that promise to substantially improve our current performance gains, and could be implemented in a production-ready version of InterpMC.

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