

A Collection of Numerical Recipes Useful For Building Scalable Psychometric Applications

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Abstract

Next generation psychometrics will inevitably leverage advancements made available with modern computing to create psychometric innovations. Realizing this future requires that psychometricians have the full suite of skills needed to create and deploy new psychometric applications at scale. Unfortunately, psychometrics, as well as other quantitative fields, has not fully embraced well-known numerical recipes for computing at scale and this is a significant barrier. Commonly, textbook representations of mathematical solutions are confused to be the same computational strategy for implementation or trivial approximations are often used when faced with complex scenarios, such as simulation or basic integration techniques. Overcoming these limitations will require some focused conversations envisioning an advanced skill set for psychometricians involving numerical methods, psychometric computing, and software development. This work sets out to focus on only one of those components and a collection of numerical recipes for common psychometric problems is provided.

Keywords: computational psychometrics; item response theory; gaussian quadrature; linear models; numerical analysis; psychometric data scientist

Introduction

Psychometrics is evolving and increasingly becoming a complex and multidisciplinary field. The ideas and potential in front of next generation psychometricians presents a future for them to leverage advances in modern computing to build, scale, and implement redesigned assessment models that better align with how individuals learn in modern society and demonstrate what they have learned in new ways (von Davier, Deonovic, Yudelson, Polyak, & Woo, 2019).

Stepping towards these advancements will inevitably require a concentrated focus on building a new type of psychometric innovator. This new innovator is of course trained in traditional measurement topics, but also has significant computational acumen coupled with the full suite of skills

needed to build and deploy software applications to make those ideas accessible for broad audiences. Consider, as an example, one computational psychometrics field guide proposed by ACT (ACTNext, 2019). The ideas in this guide have greater emphasis on topics that seemingly resemble pure computer science than topics currently emphasized in traditional measurement programs.

The future of psychometrics will be driven by the modern advantages provided by new technologies in at least three ways. First, some routine psychometric tasks can be reformulated to capture efficiencies so that the computing time is significantly reduced (Lee, Leemaqz, & McLachlan, 2016) or provide avenues to directly tackle complex problems with deterministic approaches rather than reverting to stochastic approximations. Second, psychometricians sufficiently familiar with computational methods and software development will lead the advancement with new ideas and innovations for learning and assessing that will be realized as the new psychometrics. Third, the development of psychometric applications now can be more easily distributed at scale using web-based applications allowing new ideas to be more widely accessible.

One possible negative side effect of modern computing is that it is almost too easy to rely on auto-scaling in cloud-based infrastructures or distributed computing concepts to split up tasks. These methods allow for a developer to easily scale up any computational task by quickly involving shared computing power, which inevitably involves greater cost. However, it is critically more important as a first step to consider the smallest and most manageable unit—the computational methods used in the underlying code to build the application. Then, good computational methods can be more readily scaled than expecting poor methods to be remedied by more powerful computing.

Numerical methods are the core at which the ideas are implemented and psychometric innovators should have a portfolio of routine methods at the ready to support the development of new ideas. Unfortunately, efficient numerical methods are not widely embraced within the field of psychometrics and, perhaps, this is rooted in how we learn. One general observation is that textbooks commonly represent psychometric solutions with scalar-based expressions or, when matrices are displayed, they are typically used to render solutions to the problem in ways that are not computationally efficient. Those representations are valuable for developing a conceptual understanding of the psychometric problem and are intended to communicate the idea of the model or its parameterization. However, those same representations are not intended to be used for computational implementation, although it appears they often are.

Bates (2004) noted that the algebraic textbook solution for the least squares model is commonly used within R code and provided details for better computational implementation. Genz and Kass (1997) suggested that statisticians often ignore superior methods for high dimensional integration, a problem also described by Antal and Oranje (2007) in psychometric software applications finding that trivial rectangular rules are common even in the National Assessment of Educational Progress (Allen, Donoghue, & Schoeps, 2001) and commercially available psychometric software.

There are, however, many interesting conversations among psychometric innovators regarding the challenges in building scalable applications. Cai, Yang, and Hansen (2011) openly discuss the specific challenges between Gauss Hermite and rectangular rules as do the developers of the R package Dire (Doran, Bailey, Buehler, & joo Lee, 2021) for estimating the marginal likelihood regression (Mislevy, 1984) which argues for a computational trade-off that may be realized by the simpler numerical integration rule relative to using a more sophisticated approach. Rosseel (2021) provides an impressive example that translates a complex problem into a feasible computational solution. This work makes good use of a standard inversion lemma to reduce the overhead associ-

ated with inverting large matrices and takes advantage of certain special structures and reduces the problem to a yield a simpler solution and is an exemplar of psychometric computing.

One challenge is that these conversations are typically focused on a specific application or problem. In contrast, the work presented here is intentionally problem agnostic with the goal of organizing a general class of useful solutions for common psychometric problems to be used as templates for creating new solutions. For this reason, this work is more focused on application and is therefore an organization of tested numerical approaches with the hope that they become more familiar and applied in psychometric computing. Here the concept of computational psychometrics is explored through the lens of integrating psychological measurement with ways to computationally implement, showcase, and advance new ideas into better ways to assess learning and cognition. This work does not set out to formally prove theorems or engage in deriving estimators and proving their properties other than one simple proof in the appendix.

In many respects, numerical analysis techniques are situational and implemented as preference by the innovator suitable for a given condition. For instance, the examples described here make use of a certain decomposition; however, others may have legitimate reasons for others. Hence, there is no motivation to oversell these specific concepts as the most desirable for any scenario. In fact, that’s exactly the opposite point. Instead, the underlying motivation is to encourage staring at a problem and finding a recipe through exploration and testing that yields a stable and reduceable structure to implement a complex procedure at scale. In other words, the idea proposed is that psychometricians need a toolkit allowing them to make a big problem small and the ideas presented here are generally useful, but not intended as the most ideal for every scenario.

Anecdotally, the work described here is a reflection of many conversations over the years answering the basic questions, “how do I write code for this?” or “why is my code taking hours (or days) to reach a stationary point?” For this reason, the hope is to broaden those conversations to a larger audience and provide what is hoped to be a useful resource for those interested in computing complex psychometric models. The recipes provided here are introduced in the framework of familiar and simple problems and the ideas are advanced sequentially to more complex scenarios. The work ends with two examples of methods not traditionally found in standard software simply to illustrate implementation of computational ideas and not to center attention on those specific methods.

Notation and Terminology

Standard notation is generally used throughout and the terms and notation are defined within each topic. For instance, matrices assume the general notation of uppercase bold, \mathbf{A} , determinants of matrices, $|\cdot|$, vectors are generally lowercase bold, \mathbf{x} , elements within sets $\{x_1, x_2, \dots, x_n\}$, and functions of variables with a general form $f(\cdot)$.

Distinctions in terminology are made throughout using “algebraic” representation and “computational” representation to mean very different things. Specifically, “algebraic solution” is used to mean the way a solution to a problem is written. For instance, the system of linear equations generically written as $\mathbf{Ax} = \mathbf{b}$ is commonly thought of having the algebraic solution $\mathbf{x} = \mathbf{A}^{-1}\mathbf{b}$. Instead, a “computational solution” is a numerical recipe that provides a stable way to implement a solution in software to estimate parameters that may not resemble the algebraic solution.

Introductory Concepts For Dealing With Correlated Data

Whitening Transformations for Correlated Variables

A concept first introduced and used throughout almost all subsequent recipes is referred to as the whitening transformation. Whitening is a process of transforming a collection of correlated vectors into a collection of uncorrelated vectors (Kessy, Lewin, & Strimmer, 2018) with the aid of a suitable decomposition. The process is critical to the solution of linear systems and for fast quadrature routines and is generally helpful for reducing complex problems into more computationally feasible solutions.

The concept is introduced as follows. Let \mathbf{Y} be a matrix of correlated variables from a P -dimensional multivariate normal, $\mathbf{Y} \sim \mathcal{N}_p(\boldsymbol{\mu}, \boldsymbol{\Sigma})$ where $\boldsymbol{\Sigma}$ is symmetric and positive definite. Now find a decomposition that satisfies $\boldsymbol{\Sigma} = \mathbf{L}\mathbf{L}'$ where \mathbf{L} is lower triangular known as the Cholesky factor and \mathbf{L}' is its transpose. This provides that the change of variable transformation can be used as $\mathbf{X} = \mathbf{Y}\mathbf{L}'^{-1}$ so that the variables in the matrix \mathbf{X} are “whitened” or “de-correlated”. Note that \mathbf{L}'^{-1} is used as the algebraic representation; finding the inverse of a triangular matrix is simple using forwards or backwards substitution. Notice this can also be used to create correlated vectors when the operation is reversed, $\mathbf{X}\mathbf{L}' = \mathbf{Y}$, which may be useful for data generation, simulation, or applications such as high dimensional integration.

There are many useful decompositions in numerical analysis (Searle, 1982; Zhang, 1999). However, the Cholesky is broadly helpful to speed computations and some work has demonstrated its potential advantage over others, such as the QR (Lira, Iyer, Trindade, & Howle, 2016) when used for solving normal equations. Some applications have fully adopted a sparse Cholesky decomposition (Davis & Hager, 2005) for linear mixed models (Bates, 2004b) with tremendous success. For these reasons, the Cholesky is the primary decomposition used throughout.

Multivariate Densities of Whitened Data

The multivariate normal density function has an important role in psychometric practice. It is possible to reduce this from a complex problem into separable parts and show how the multivariate density can be computed as a product of transformed univariate densities. This change of variable transformation has broader applications for psychometric computing in maximization methods or dealing with high-dimensional integration problems.

The general form of the P -dimensional multivariate normal for a non-diagonal covariance matrix $\boldsymbol{\Sigma}$ is

$$f_y(y_1, \dots, y_p) = \frac{\exp(-\frac{1}{2}(\mathbf{y} - \boldsymbol{\mu})'\boldsymbol{\Sigma}^{-1}(\mathbf{y} - \boldsymbol{\mu}))}{\sqrt{(2\pi)^p |\boldsymbol{\Sigma}|}} \quad (1)$$

If the mean vector is $\boldsymbol{\mu} = \mathbf{0}$ and $\boldsymbol{\Sigma} = \text{diag}(1, \dots, 1)$, then the density of the multivariate normal is equal to the product of the marginal densities. Let φ denote the standard normal probability density function (p.d.f.), then

$$f_y(y_1, \dots, y_p) = \prod_{i=1}^p \varphi(y_i) \quad (2)$$

Note that (1) and (2) are related only under the special case of uncorrelated variables with unit variance. With correlated variables, rewrite (1) as

$$f_y(y_1, \dots, y_p) = \frac{\exp(-\frac{1}{2}(\mathbf{y} - \boldsymbol{\mu})' \mathbf{L}^{-1} \mathbf{L}'^{-1} (\mathbf{y} - \boldsymbol{\mu}))}{\sqrt{(2\pi)^p |\mathbf{L} \mathbf{L}'|}}. \quad (3)$$

Now let $\mathbf{x}^{*'} = \mathbf{y}' \mathbf{L}^{-1}$ such that the elements of $\mathbf{x}^{*'} = \{x_1^*, \dots, x_p^*\}$, then

$$f_y(y_1, \dots, y_p) = \frac{\exp(-\frac{1}{2} \mathbf{x}^{*'} \mathbf{x}^*)}{\sqrt{(2\pi)^p |\mathbf{L}^2|}} \quad (4)$$

$$= |\mathbf{L}|^{-1} \left(\prod_{i=1}^p \varphi(x_i^*) \right) \quad (5)$$

This transform expresses (5) into the form of (2), thus showing how the multivariate density can be computed as a product of the (transformed) marginal densities.

Computational Concepts for Linear Systems

This section explores efficient computational recipes for linear models to create simple, reducible structures to build scalable psychometric algorithms. It's best to begin with a simple, common problem—ordinary least squares (OLS), that will serve as a foundation and subsequently show how more complex problems can be reduced to the same OLS solution. While the examples are provided within the context of linear regressions, the same concepts are used widely across psychometric practice. For example, these same methods are often used in the expectation-maximization (EM) algorithm during the M-step, in natural language processing problems using latent semantic analysis, solutions to factor analysis problems, in highly parameterized item calibrations problems using the newton-raphson algorithm, and within iteratively reweighted least solutions for generalized linear models. These are just a subset of the possible applications.

Computing Ordinary Least Squares

The standard representation of the linear least squares problem is via the following normal equations (McCulloch & Searle, 2001)

$$\mathbf{X}' \mathbf{X} \boldsymbol{\beta} = \mathbf{X}' \mathbf{y} \quad (6)$$

where \mathbf{X} is an $n \times p$ design matrix of full column rank, $\boldsymbol{\beta}$ is a $p \times 1$ vector of unknown parameters to be estimated, and \mathbf{y} is an $n \times 1$ vector of observed outcomes. Because $\mathbf{X}' \mathbf{X}$ is $p \times p$, it implies it is invertible and the following algebraic solution typically follows in texts

$$\hat{\boldsymbol{\beta}} = (\mathbf{X}' \mathbf{X})^{-1} \mathbf{X}' \mathbf{y}. \quad (7)$$

The algebraic representation of the solution for the estimates of the fixed effects in Equation (7) is generally conflated with its computational solution. It is in fact unnecessary to actually invert

$X'X$; instead it is more reliable to find a decomposition for $X'X$ and then find the parameter estimates via triangular systems rather than using matrix inversion.

One stable approach is to find the decomposition, $X'X = LL'$, and then rewrite (6) as $LL'\beta = X'y$. This expression provides a simple solution via triangular systems requiring that we first forward solve for z and then backsolve for β as

$$Lz = X'y, L'\beta = z. \quad (8)$$

Obtaining least squares estimates is very simple. However, it serves as a baseline for computing and as subsequently shown, when more complex models can be rewritten into the simplified form of (6), then they also inherit the simple solution shown in (8).

Computing Generalized Least Squares Using Decompositions

The generalized least squares (GLS) problem is an incrementally more advanced model than its OLS counterpart. The standard textbook normal equations for the GLS model is

$$X'\Sigma^{-1}X\beta = X'\Sigma^{-1}y \quad (9)$$

where Σ is a symmetric ($n \times n$) positive definite covariance matrix and all others matrices are as previously noted. The algebraic solution for the GLS approach is commonly shown to involve matrix inversions as follows

$$\hat{\beta} = (X'\Sigma^{-1}X)^{-1}X'\Sigma^{-1}y. \quad (10)$$

This algebraic representation is conceptually useful, but does not make for an inefficient computational implementation. In particular, the solution to (10) implies inverting the matrix Σ and then again inverting $(X'\Sigma^{-1}X)$. The dimensions of Σ are very large and the GLS problem often involves an iterative solution with elements of Σ changing over the iterations.

The concept of whitening can again be used and provides that the data can be “de-correlated” via the transformation $X^* = XL'^{-1}$ where $\Sigma = LL'$. This provides a way to rewrite the normal equations as

$$X'L^{-1}L'^{-1}X\beta = X'L^{-1}L'^{-1}y \quad (11)$$

$$X^{*'}X^*\beta = X^{*'}y^* \quad (12)$$

where $X^* = XL'^{-1}$ and $y^* = yL'^{-1}$. This whitening transform writes (12) into the form (6) so that it can inherit the same solution shown in (8).

If (12) were used in an iterative process and elements of Σ were updated at each step, then recomputing its Cholesky factor repeatedly would layer additional expense. This is where numerical recipes become situational and we encourage staring at Σ to determine if it can be separated and reduced into a simple structure. We provide one such example in the section on linear mixed models where a very large matrix is repeatedly updated in an iterative algorithm and finding its inverse at each iteration is costly. Instead, a standard inversion lemma is used and it is extended to exploit some special characteristics of that matrix allowing a complicated problem to be reduced into a very trivial computation implementation.

Computing Weighted Least Squares

The weighted least squares (WLS) problem is a special case of the GLS problem when the matrix Σ in the preceding example is diagonal. In this case, it is entirely unnecessary to actually create and store the entire matrix Σ and instead only its diagonal elements, v_{jj} , are needed. In this case there is a special matrix property providing that $\mathbf{X}'\Sigma^{-1}\mathbf{X} = \mathbf{X}'\mathbf{X}^*$ where \mathbf{X}^* is arrived at \mathbf{X}^* via the elementwise calculation of $1/v_{jj}$ over each column of the matrix \mathbf{X} .

Performing this elementwise calculation to the matrix \mathbf{X} on the left and right hand side of Equation 9 when Σ is diagonal yields

$$\mathbf{X}'\mathbf{X}^*\boldsymbol{\beta} = \mathbf{X}'^*\mathbf{y}. \quad (13)$$

Equation (13) also is now written in the form of (6) and inherits the same computational solution.

Linear Mixed Models with Henderson's Method

The linear mixed model is generally written with the form (Laird & Ware, 1982)

$$\mathbf{Y} = \mathbf{X}\boldsymbol{\beta} + \mathbf{Z}\boldsymbol{\theta} + \epsilon \quad (14)$$

where \mathbf{Y} is an $n \times 1$ vector of outcomes, \mathbf{X} is an $n \times p$ model matrix for the fixed effects, $\boldsymbol{\beta}$ is a $p \times 1$ vector of estimates for the fixed effects, $\mathbf{Z} = [\mathbf{Z}_1, \mathbf{Z}_2, \dots, \mathbf{Z}_Q]$ is an $n \times Q$ model matrix for the random effects, $\boldsymbol{\theta}' = [\boldsymbol{\theta}'_1, \boldsymbol{\theta}'_2, \dots, \boldsymbol{\theta}'_Q]$ is the vector of random effects, and ϵ as the residual error term. The normal equations for the linear mixed model is (McLean, Sanders, & Stroup, 1991)

$$\begin{bmatrix} \mathbf{X}'\Omega^{-1}\mathbf{X} & \mathbf{X}'\Omega^{-1}\mathbf{Z} \\ \mathbf{Z}'\Omega^{-1}\mathbf{X} & \mathbf{Z}'\Omega^{-1}\mathbf{Z} + \mathbf{G}^{-1} \end{bmatrix} \begin{bmatrix} \boldsymbol{\beta} \\ \boldsymbol{\theta} \end{bmatrix} = \begin{bmatrix} \mathbf{X}'\Omega^{-1}\mathbf{y} \\ \mathbf{Z}'\Omega^{-1}\mathbf{y} \end{bmatrix}. \quad (15)$$

The linear mixed model iterates to yield a final solution with the following general steps:

1. Construct starting values for the variances of the random effects including σ_ϵ^2 and σ_q^2 for all levels of q . These are used as elements in the diagonal matrices Ω and \mathbf{G} , respectively.
2. Solve the linear system for $\boldsymbol{\beta}$ and $\boldsymbol{\theta}$
3. Update the values of the variances of the random effects including σ_ϵ^2 and σ_q^2
4. Iterate between steps 2 and 3 until the difference between any of the parameter estimates is less than a prespecified convergence criterion.

Given the representation of Equation (15) there is a temptation to assume a naive solution to the linear system using matrix inversion

$$\begin{bmatrix} \hat{\boldsymbol{\beta}} \\ \hat{\boldsymbol{\theta}} \end{bmatrix} = \begin{bmatrix} \mathbf{X}'\Omega^{-1}\mathbf{X} & \mathbf{X}'\Omega^{-1}\mathbf{Z} \\ \mathbf{Z}'\Omega^{-1}\mathbf{X} & \mathbf{Z}'\Omega^{-1}\mathbf{Z} + \mathbf{G}^{-1} \end{bmatrix}^{-1} \begin{bmatrix} \mathbf{X}'\Omega^{-1}\mathbf{y} \\ \mathbf{Z}'\Omega^{-1}\mathbf{y} \end{bmatrix}.$$

As in the linear models discussed so far, the Cholesky factor of the leftmost matrix in (15) can be obtained

$$LL' = \begin{bmatrix} X'\Omega^{-1}X & X'\Omega^{-1}Z \\ Z'\Omega^{-1}X & Z'\Omega^{-1}Z + G^{-1} \end{bmatrix} \quad (16)$$

and let $\Theta' = [\beta' \ \theta']$ and

$$y^* = \begin{bmatrix} X'\Omega^{-1}y \\ Z'\Omega^{-1}y \end{bmatrix}.$$

This is now written in the general form used in the OLS description and shares the same solution using (8). First solve $Lz = y^*$ for z and then solve $L'\Theta = z$ for Θ . Of course, this is only one iteration and solutions for the mixed model iterate with updated values for the variance components. Note, that when all elements of the matrix $G = 0$, this reduces to the partitioned fixed effects model.

Variance Estimation with Woodbury Identity and Decompositions for Matrix Inversions

Because Henderson's method is an iterative procedure, values for the variance components are revised at each step. This can be accomplished via the following, which is computed at each iteration of the algorithm (McCulloch & Searle, 2001)

$$T = \left[I + \left(Z'Z - Z'X(X'X)^{-1}X'Z \right) D \right]^{-1} \quad (17)$$

where I is an identity matrix with dimensions equal to the model matrix for the random effects and $D = \sigma_{e,i-1}^{-2}G$ and all other matrices are as above. The residual variance at the i th iteration is

$$\sigma_{e,i-1}^2 = \frac{y'e}{N - p} \quad (18)$$

and the estimate of the variance at level q at the i th iteration is:

$$\sigma_q^2 = \frac{\theta_q^{i-1'} \theta_q^{i-1}}{m_q - tr(T_q)} \quad (19)$$

where θ_q are the predictions of the random effects at level q , m_q are the number of units in level q , and $tr(T_q)$ is the trace of the matrix T_q , which is the block of the matrix T corresponding to the q th block.

The challenge here is that the square matrix T can be very large. As a result, it is extremely expensive to find its inverse inside an iterative problem, thus rendering this particular approach computationally prohibitive unless a reduceable structure can be identified. In some cases, there is an alternative method for taking the inverse of a matrix with this structure via the Woodbury Matrix Identity (Woodbury, 1950) that provides a very efficient route for computation. First, Equation 17 is rewritten as:

$$T = \left(I + Z'ZD - Z'X(X'X)^{-1}X'ZD \right)^{-1} \quad (20)$$

where D is a diagonal matrix in the case of random intercepts and within each block q is constant along the diagonal. The Woodbury Identity provides that T can be computed as:

$$\mathbf{T} = \mathbf{A}^{-1} + \mathbf{A}^{-1}\mathbf{B}(\mathbf{P} - \mathbf{B}'\mathbf{D}\mathbf{A}^{-1}\mathbf{B})^{-1}\mathbf{B}'\mathbf{D}\mathbf{A}^{-1} \quad (21)$$

where $\mathbf{A} = \mathbf{I} + \mathbf{Z}'\mathbf{Z}\mathbf{D}$, $\mathbf{B} = \mathbf{Z}'\mathbf{X}$, and $\mathbf{P} = \mathbf{X}'\mathbf{X}$. This identity provides that the inverse of \mathbf{T} can be computed by separately finding the inverse of \mathbf{A} and then also of $\mathbf{P} - \mathbf{B}'\mathbf{D}\mathbf{A}^{-1}\mathbf{B}$. The latter of the two has very small dimensions, only $p \times p$ where p is the number of fixed effects in the model. The portion $\mathbf{I} + \mathbf{Z}'\mathbf{Z}\mathbf{D}$ is a bit more cumbersome because the dimensions of the model matrix for the random effects is $Q \times Q$, where Q is the number of columns in the model matrix for the random effects which can be extremely large.

The components \mathbf{I} and $\mathbf{Z}'\mathbf{Z}$ are fixed and never change over the iterative process, only \mathbf{D} is changing at each iteration. For nested random effects, $\mathbf{Z}'\mathbf{Z}$ is block diagonal but it has no special structure when the random effects are fully or partially crossed. This fact is motivation to explore whether special structures exist that might be exploited in order to more easily find \mathbf{A}^{-1} given that it will be performed within each iteration. Of course, it is a well-known result that the sum of inverses is not equal to the inverse of the sum. However, both \mathbf{I} and \mathbf{D} have special structures to exploit—both are diagonal within a block of q for nested random effects and have constant values along the diagonal.

These special structures make it possible to find \mathbf{A}^{-1} in a trivial way for some variations of the linear mixed model. When dealing with nested random effects, $(\mathbf{Z}'\mathbf{Z})^{-1}$ is block-diagonal and its inverse can be found by taking the inverse one block at a time. Within each block, the diagonal elements of \mathbf{D} are constant. It's pointed out here that the following works in the special case of nested random effects for random intercepts and is not globally generalizable (see appendix). However, the emphasis in this work is to identify special structures for a given scenario and exploit them to reduce computational burden. The inverse of the sum provided here may also have broader applications in other types of problems and so it is a useful exercise to document the condition.

For convenience, the subscript on \mathbf{Z} is dropped, but note that this performed on each block, or perhaps on large \mathbf{Z} , when it too has a special structure. First, precompute and store the eigendecomposition for $\mathbf{Z}'\mathbf{Z}$, which is the most expensive part of the calculation. Then, all subsequent iterations treat the eigenvectors as fixed and are reused and the eigenvalues are updated repeatedly with virtually no computational overhead. The inverse of the sum is then simply the inverse of the eigenvalues as

$$(\mathbf{Z}'\mathbf{Z}\mathbf{D} + \mathbf{I})^{-1} = \mathbf{A}^{-1} = \mathbf{Q}_1\boldsymbol{\lambda}^{*-1}\mathbf{Q}_1' \quad (22)$$

where $\mathbf{Z}'\mathbf{Z} = \mathbf{Q}_1\boldsymbol{\lambda}_1\mathbf{Q}_1'$ and $\boldsymbol{\lambda}^* = \text{diag}\{\lambda_{11}d + 1, \lambda_{12}d + 1, \dots, \lambda_{n1}d + 1\}$. Proof for (22) is provided in the appendix. Now \mathbf{A}^{-1} is obtained in a much less expensive way by reusing, \mathbf{Q}_1 , and simply updating $\boldsymbol{\lambda}^*$ at each iteration, which is a nominal task given that it is diagonal. So, pulling together the Woodbury Identity and the decomposition derived for Equation (22), the inverse of the very big matrix \mathbf{T} is easily computed over and over again with little overhead.

Scalable Methods for Integration

Psychometricians routinely encounter challenging marginal likelihood problems or posterior distributions that require numerical integration methods. These methods approximate difficult integrals by summing over a finite number of points. Integrals with a gaussian factor are rather common in

psychometric applications and, as such, the Gauss Hermite rule may be used for integrals of the following form over the domain $(-\infty, \infty)$

$$\int f(\theta)e^{-\theta^2}d\theta \approx \sum_{q=1}^Q f(\theta_q)w_q. \quad (23)$$

where summation is over $\mathcal{P} = \{\theta_q, w_q\}_{q=1}^Q$, the so called nodes, θ_q , and weights, w_q , for $q = 1, 2, \dots, Q$. The choice of nodes and weights in \mathcal{P} is optimal under the Gauss Hermite rule and will generally outperform other methods (Quarteroni, Saleri, & Gervasio, 2010). Unfortunately, psychometric practice seems to have codified the basic rectangular rule with fixed, equally spaced points from the domain of $(-4, 4)$ (Bock & Mislevy, 1982) despite its serious limitations in most cases for at least two reasons. First, the equally spaced points may not be sampled in the best region of the integrand and second it lacks the computational advantage of precision until an extremely large number of fixed points are used. Building scalable applications requires approaches with minimal computational overhead that achieve precision rapidly.

One Dimensional Gaussian Quadrature

Conveniently, psychometricians encounter scenarios where the weight function is a standard normal density

$$\int f(\theta) \frac{e^{-\theta^2/2}}{\sqrt{2\pi}} d\theta \approx \pi^{-1/2} \sum_{q=1}^Q f(\sqrt{2}\mathbf{L}\theta_q + \tilde{\theta})w_q \quad (24)$$

where $\tilde{\theta}$ is a location parameter for centering (Liu & Pierce, 1994)¹. The approximation in (24) implements a transformation on θ from (23) and adapts to the best region of the integrand—hence it is the adaptive Gauss Hermite rule. The expression in (24) suggests $\tilde{\theta}$ and values to construct $\Sigma = \mathbf{L}\mathbf{L}'$ are known, or at least approximately known, yet those are the latent quantities we are trying to estimate. We return to an approach to consider for replacing these with provisional values in a later section.

Multivariate Gaussian Quadrature

The one dimensional Gauss Hermite rule can be extended to the case for integrals of the general form in multiple dimensions (Chowdhary, Salloum, Debusschere, & Larson, 2015; Jäckel, 2005; Judd, Maliar, & Maliar, 2011; Stringer, 2021)

$$\int \dots \int f(\boldsymbol{\theta})\psi(\boldsymbol{\theta})d\boldsymbol{\theta}. \quad (25)$$

where $\boldsymbol{\theta} = \{\theta_1, \dots, \theta_p\}'$ and continue with the assumption $\psi(\boldsymbol{\theta})$ is multivariate normal. The integral in (25) is often intractable and requires approximation. This may generally be approximated with nodes rotated to reflect a covariance structure and possibly shifted over a mean. That rotation

¹The \mathbf{L} in (24) is actually a scalar, σ , but is left as a matrix so it appears as a special case of (26).

can be implemented with the Cholesky factor of a covariance matrix as (Jäckel, 2005; Judd et al., 2011; Stringer, 2021)

$$\int \dots \int f(\boldsymbol{\theta}) \psi(\boldsymbol{\theta}) d\boldsymbol{\theta} \approx \pi^{-P/2} \sum_{q_1=1}^Q \dots \sum_{q_p=1}^Q f(\sqrt{2}\mathbf{L}(\theta_{q_1}, \dots, \theta_{q_p})' + (\tilde{\theta}_{q_1}, \dots, \tilde{\theta}_{q_p})') w_q^* \quad (26)$$

where θ_{qp} is node $q = 1, 2, \dots, Q$ in dimension $p = 1, 2, \dots, P$, w_{qp} is the corresponding weight, $w_q^* = \prod_{p=1}^P w_{qp}$, and $\tilde{\theta}_{qp}$ is used to mean some points around which the nodes are centered. The nodes are expansions of the points in the one-dimensional set, \mathcal{P} , via a cartesian product rule to form a multivariate grid, \mathbf{G} , with dimensions $Q^P \times P$. Hence, the q th row of \mathbf{G} is $\mathbf{g}_q = (\theta_{q1}, \dots, \theta_{qP})$. Note that (24) is simply a special case of (26), hence no new concepts are needed for implementation beyond what has been described for the one-dimensional scenario.

The extremely large number of quadrature points is what renders this approach challenging as, hueristically, this is effectively a two step process. Step 1 requires evaluating the function $f(\cdot)$ over all Q^P nodes and then step 2 is the summation of (26) making use of those values over all rows of the grid \mathbf{G} . Even with a modest selection for Q , the exponential growth in the function evaluations quickly becomes large with increasing P . Hence, we might explore ways in which we evaluate $f(\cdot)$ as few times as possible and whether fewer function evaluations than $Q^P \times P$ can be used.

Applications to Psychometric Problems

The previously discussed methods are computational templates that can be used to build and scale psychometric applications. Two examples are provided here with supporting code in the appendix to bring some life to those concepts. The intent is not to center attention on these specific examples; rather these examples are used because they are approaches that typically require computing something complex and are not routinely found in standard software. For this reason, they are attractive for exploring computationally efficient approaches.

Example 1: Estimating the Error in Variables Linear Model

Estimates of examinee performance from a psychometric instrument are noisy reflections of a true, latent trait. The observed measures are subsequently used in various ways and often used as inputs in regression models of some form. When using these observed estimates, the assumptions of the Gauss-Markov theorem do not hold and the parameter estimates are then inconsistent unless corrections are made (Doran, 2014; Lockwood & McCaffrey, 2020; Nab, van Smeden, Keogh, & Groenwold, 2021).

These models are chosen as an example because with different assumptions about the nature of the measurement error (e.g., homoscedastic versus heteroscedastic), the models may take different forms and custom software is often written to obtain parameter estimates. One such example is the `mecor` (Nab et al., 2021) package in R and a code review of its implementation shows that the algebraic representation and computational representation are one in the same.

One form of the linear EiV regression is commonly shown to be (StataCorp, n.d.)

$$(\mathbf{X}'\mathbf{X} - \mathbf{S})\boldsymbol{\beta} = \mathbf{X}\mathbf{y} \quad (27)$$

where, in the case of known estimates for the reliability, the matrix $\mathbf{S} = \text{diag}(N(1-r_1)\sigma_1^2, \dots, N(1-r_j)\sigma_j^2)$ where r_j is the reliability of the j th variable, N is the number of individuals, and σ_j^2 is the sample variance of the j th variable in the model matrix \mathbf{X} . Now, find $\mathbf{X}'\mathbf{X} - \mathbf{S} = \mathbf{L}\mathbf{L}'$ and obtain estimates using (8). Code to implement a simple error-in-variables model with known reliability is provided in the appendix as an example. It's useful to note that (27) can be extended for the linear mixed model and the computational approach previously shown can be applied.

Example 2: Estimating a High-Dimensional EAP

Suppose we are interested in estimating the ability of an examinee using a P -dimensional EAP estimator. This is chosen as an example because it is a desirable estimator that continues to be challenging for psychometricians to implement (Chalmers, 2012; Ferrando & Lorenzo-Seva, 2016). Let the general form of the integral be represented as

$$\hat{\boldsymbol{\theta}} = \frac{\int \cdots \int \boldsymbol{\theta} L(\mathbf{z}|\boldsymbol{\theta}) \psi(\boldsymbol{\theta}; \boldsymbol{\mu}, \boldsymbol{\Sigma}) d\boldsymbol{\theta}}{\int \cdots \int L(\mathbf{z}|\boldsymbol{\theta}) \psi(\boldsymbol{\theta}; \boldsymbol{\mu}, \boldsymbol{\Sigma}) d\boldsymbol{\theta}} \quad (28)$$

where $\psi(\boldsymbol{\theta}; \boldsymbol{\mu}, \boldsymbol{\Sigma})$ is the multivariate normal density function and the likelihood function (see appendix) is a product of domain-specific likelihoods as

$$L(\mathbf{z}|\boldsymbol{\theta}) = L(\mathbf{z}_1|\theta_1)L(\mathbf{z}_2|\theta_2), \dots, L(\mathbf{z}_p|\theta_p).$$

We assume individuals have some idiosyncratic ability contributing to their response on each dimension and that some test items are uniquely associated with dimension 1 (θ_1), others with dimension 2 (θ_2), and so on for the P th dimension, $\boldsymbol{\theta} = (\theta_1, \theta_2, \dots, \theta_P)'$.

The likelihood function and its corresponding population distribution in (28) are not in the same parametric family, as such the integral cannot be evaluated analytically. Here three options are explored that make for possible implementation that vary in the number of times the function $L(\mathbf{z}|\boldsymbol{\theta})$ is computed. Once the likelihood is evaluated, then step 2 is effectively direct implementation of the sum over all rows of \mathbf{G} . Sparse grids or “pruning” some values are options for reducing the number of points over which the function is evaluated, but those are not explored here.

The first canonical option is to use (26) directly. This involves using the Cholesky factor to adjust the nodes to absorb the population covariance structure and then, possibly shift the nodes over a new centered point. This works quite well, and in fact is implemented in the code in the appendix. This however, has one downside worth noting that makes this computationally somewhat less attractive.

Initially (before rotating with the Cholesky factor), the nodes in \mathbf{G} are patterned such that marginally (i.e., each column in \mathbf{G}) the Q unique nodes are repeated over and over. When the rows of this matrix are rotated via $\mathbf{L}(\theta_{q1}, \dots, \theta_{qp})'$, then the uniqueness is lost exponentially over each column. That is, column 1 has Q unique nodes repeated, column 2 has only Q^2 unique nodes, column 3 would have Q^3 unique nodes and so on up to the P th column.

This reduced pattern of unique nodes creates a larger burden in that we must compute the values of the likelihood at each node and that IRT likelihood is somewhat expensive as it involves computing exponents. In column 1, it is feasible to compute the value of the likelihood at the Q

unique points and then simply copy those values into the other positions in column 1 where the same value of the node appears. That is, we would compute the likelihood only Q times and then recycle those values by copying them into other positions where the same node appears in the grid. In column 2, we must compute the likelihood at all Q^2 unique nodes and then the values could then be copied into the positions of this column where the same node appears again. The upside here is that the method yields a good approximation of the integral. The downside is the computational expense is very large. In this case, we would evaluate the likelihood $Q + Q^2 + Q^3, \dots, Q^P$ times.

A second option for computing this integral could be to use a stochastic quadrature routine. In this process the nodes are random draws from $\mathcal{N}_p(\mathbf{M}, \mathbf{I}_P)$, where \mathbf{I}_P is a P -dimensional identity matrix again creating the grid \mathbf{G} , but in this instance it would have dimensions $R \times P$ where R is the number of random variates chosen and it is not an exponential function of R . In this manner, the variates in \mathbf{G} would reflect the population covariance after being premultiplied by the Cholesky factor (see introductory section) when the summation in (26) is used to evaluate the integral. This stochastic approach makes the computation reasonable and its precision is highly dependent on the law of large numbers such that $\hat{\theta}_p \xrightarrow{a.s.} \theta_p$ as $R \rightarrow \infty^2$. In this case, we would evaluate the likelihood $R \times P$ times.

Perhaps a third and seemingly very rapid approach can be considered. First, rewrite the integral within (28)

$$\int \dots \int L(\mathbf{z}|\boldsymbol{\theta})\psi(\boldsymbol{\theta}; \boldsymbol{\mu}, \boldsymbol{\Sigma})d\boldsymbol{\theta} = \int \dots \int \frac{L(\mathbf{z}|\boldsymbol{\theta})\psi(\boldsymbol{\theta}; \boldsymbol{\mu}, \boldsymbol{\Sigma})}{\tilde{\psi}(\boldsymbol{\theta}; \tilde{\boldsymbol{\mu}}, \tilde{\boldsymbol{\Sigma}})}\tilde{\psi}(\boldsymbol{\theta}; \tilde{\boldsymbol{\mu}}, \tilde{\boldsymbol{\Sigma}})d\boldsymbol{\theta} \quad (29)$$

which has approximation

$$\hat{\theta}_p = \frac{\sum_{q=1,1}^Q \dots \sum_{q=p,1}^Q \theta_{p,q} L(\mathbf{z}_1|\theta_{1,q}) L(\mathbf{z}_2|\theta_{2,q}) \dots L(\mathbf{z}_p|\theta_{p,q}) \phi(\boldsymbol{\theta}_q; \boldsymbol{\mu}', \boldsymbol{\Sigma}') w_q^*}{\sum_{q=1}^Q \dots \sum_{q=1}^Q L(\mathbf{z}_1|\theta_{1,q}) L(\mathbf{z}_2|\theta_{2,q}) \dots L(\mathbf{z}_p|\theta_{p,q}) \phi(\boldsymbol{\theta}_q; \boldsymbol{\mu}', \boldsymbol{\Sigma}') w_q^*} \quad (30)$$

where

$$\phi(\boldsymbol{\theta}_q; \boldsymbol{\mu}', \boldsymbol{\Sigma}') = \frac{\psi(\boldsymbol{\theta}_q; \boldsymbol{\mu}, \boldsymbol{\Sigma})}{\tilde{\psi}(\boldsymbol{\theta}_q; \tilde{\boldsymbol{\mu}}, \tilde{\boldsymbol{\Sigma}})}. \quad (31)$$

where $\tilde{\psi}(\cdot)$ is a proposal distribution. This expression resembles an importance sampling approach and has been termed importance Gauss Hermite (IGH) method (Elvira, Martino, & Closas, 2021) and other applications have used similar approaches (Antal & Oranje, 2007; Tuerlinckx, Rijmen, Verbeke, & De Boeck, 2006; Pinheiro & Bates, 1994). The value of the IGH approach is that the weights are modified to have greater importance for certain values of the integrand, in which case fewer quadrature points may be needed to evaluate the integral. The trade-off is that some calculations involving the densities in $\phi(\cdot)$ is required, though that additional cost is seemingly smaller than if more quadrature points were used and importance-based methods may be valuable for reducing computational burden (Ackerberg, 2009).

The additional benefit is that the original (unrotated) Gauss Hermite nodes are used in (30). We can exploit the uniqueness in the repeating values of the nodes in each column and compute the value of the likelihood only Q times in each dimension and then copy those values into other

²The asymptotics here also depend on the number of test items being very large and not only the number of variates used for evaluation

positions of the column where they repeat. As a result, the likelihood is evaluated only $Q \times P$ times and not $Q^P \times P$ times! Assuming $Q < R$, this approach evaluates the likelihood the smallest number of times of all three cases.

Naturally, the question of what values are used in $\tilde{\psi}(\cdot)$ needs clarification. This choice is what has the largest impact relative to the computational burden, yet unfortunately, some experimentation is needed for choosing these values. The implementation provided in the appendix uses $\tilde{\psi}(\boldsymbol{\theta}; \mathbf{0}, \mathbf{I})$ as a template.

One final consideration for reducing some computational burden in problems of this nature might be to use an iterative process as suggested by Naylor and Smith (1982). If values for the parameters of interest were known, or at least approximately known, the nodes could be recentered via the transformation $\sqrt{2}\theta_{p,q}\sigma_p + \hat{\theta}_p$ and take another pass over the integral with the transformed nodes and weights at an increased number of quadrature points for precision could be used.

R code for implementing this problem is provided in the appendix. It is a generalizable function that can evaluate any P -dimensional problem. Some additional computational benefits could easily be realized in this code. For instance, computing the likelihood values over each column in the node array is independent and consequently is embarrassingly parallel. This code version does not reflect the parallel computing as too many factors could detract from the intent of being didactic. Users could easily extend the loop within the code to be parallel, however, this does require making additional copies of the node array which alone could be memory intensive. However, almost all test scoring applications in psychometrics are an embarrassingly parallel problem as the test score for individual k does not depend on the scores for others, it depends only on the item parameters and their item responses, both of which are treated as fixed.

Summary

The ideas here are but a subset of a large array of numerical methods; however, the hope is that the work motivates next generation psychometricians to build a portfolio of techniques that can be considered for implementation when building new ideas. The intent is for the reader to see the general advantage of efficient numerical recipes that can be more broadly used in other applications to build novel psychometric solutions and to move away from the notion that algebraic representations are also to be used as computational implementations.

This work is also intended to fit into a larger conversation of the future of psychometrics. The skill set for building scalable psychometric applications means innovators need foundational numerical approaches, but also skills in software development to prototype, finalize, and deploy at scale applications to support those ideas. Commercial cloud-based infrastructures are more readily available and provide platforms to take advantage of distributed computing. Nonetheless, the first step should be to build an efficient method and then secondly scale that method with computational advantages offered.

Appendix A: Computational Formulae

Proof for Reducing $(Z'ZD + I)^{-1}$

Let $Z'Z = Q_1\lambda_1Q_1'$ and $D = Q_2\lambda_2Q_2' = \lambda_2$ given that $D = dI$, then

$$Z'ZD + I = (Q_1\lambda_1Q_1')(Q_2\lambda_2Q_2') + I \quad (32)$$

$$= (Q_1\lambda_1Q_1')\lambda_2 + I \quad (33)$$

The commutative property for matrices holds if and only if dI , hence $Q_1'\lambda_2 = \lambda_2Q_1'$. Then, $(Q_1\lambda_1Q_1')\lambda_2 + I = (Q_1\lambda_1\lambda_2Q_1') + I$. Finally, we can take advantage of the identity matrix and co-diagonalize its elements alongside the eigenvalues $\lambda_1\lambda_2 = \lambda^* = \text{diag}\{\lambda_{11}d + 1, \lambda_{12}d + 1, \dots, \lambda_{n1}d + 1\}$. Assembling all components then gives

$$Q_1\lambda_1\lambda_2Q_1' + I = Q_1\lambda^*Q_1' \quad (34)$$

Finally, the inverse is

$$(Z'ZD + I)^{-1} = Q_1\lambda^{*-1}Q_1'. \quad (35)$$

Equation (35) is useful in iterative algorithms as it precomputes an expensive component, Q_1 , initially and then it is repeatedly reused in subsequent iterations. Because updating λ^* is simple and it is diagonal, it permits a trivial inverse. Collectively, the special structures in this instance are exploited to reduce a larger computational burden into one that is much less expensive within iterative algorithms.

General Remark on the Proof

It is important to note that preceding proof applies under conditions when $D = dI$, in which case the commutative property holds between D and some other (conformable) matrix. However, when that condition is not true, we arrive at a slightly different problem. Suppose D is diagonal, but the elements along the diagonal are not constant. Then, we may write $(Z'ZD + I)^{-1} = (Q\lambda Q' + I)^{-1} = (Q(\lambda + I)^{-1}Q')$ where now $Z'ZD = Q\lambda Q'$.

This is interesting, but not entirely helpful within an iterative algorithm as D cannot be factored out of the decomposition. Hence, it requires computing $Z'ZD$ at each iteration even though only D is changing. At the current time, there is no accepted method for the inverse of matrix sums under this condition. There are some concepts that may later be considered using eigenvalue perturbations. That is, if we can first compute the decomposition $Z'Z$, then update this with the changing values of D to yield an approximation of $Q\lambda Q'$, then we might be able to efficiently find the sum of the inverse when D is diagonal, but not with constant elements. This remains a numerical challenge to explore.

Likelihood Function Details

Each individual domain likelihood, $L(z_p|\theta_P)$, is composed of a mixture of binary and polytomous test items with known item parameters. Let j index an item and then

$$L(\mathbf{z}_p|\theta_P) = L_1(\mathbf{z}_p|\theta_P)L_2(\mathbf{z}_p|\theta_P)$$

$$L_1(\mathbf{z}_p|\theta_P) = \prod_{j \in p} \left[c_j + \frac{1 - c_j}{1 + \exp[-Da_j(\theta_P - b_j)]} \right]^{x_j} \left[1 - \left(c_j + \frac{1 - c_j}{1 + \exp[-Da_j(\theta_P - b_j)]} \right) \right]^{1-x_j}$$

where $j \in p$ is used to mean there is a collection of j items uniquely associated with dimension p , c_j is the lower asymptote of the item response curve (i.e., the guessing parameter), a_j is the slope of the item response curve (i.e., the discrimination parameter), b_j is the location parameter, x_j is the observed response to the item, and D is a constant, by default fixed at 1.7. Then the items scored in multiple categories takes the form of the graded response model as

$$L_2(\mathbf{z}_p|\theta_P) = \prod_{j \in p} p(z_{jp}|\theta_P)$$

$$p(z_i|\theta) = \begin{cases} \frac{1}{1+e^{a_j D(\theta_P - b_{j1})}}, & z = 0 \\ \frac{\frac{1}{1+e^{a_j D(\theta_P - b_{j,z+1})}} - \frac{1}{1+e^{a_j D(\theta_P - b_{jz})}}}{\frac{1}{1+e^{-a_j D(\theta_P - b_{jK})}}}, & 0 < z < K \\ \frac{1}{1+e^{-a_j D(\theta_P - b_{jK})}}, & z = K \end{cases} \quad (36)$$

where b_K denotes the k th step and the other definitions used for the binary model apply here.

Appendix B: Computational Code Examples

Code Examples For Linear Systems

```
### Simple implementation of error in variables linear model
eivreg <- function(y,X,r){
  vars <- apply(X, 2, var)
  N <- length(y)
  S <- diag((1-r)*vars * N)
  XtX.S <- crossprod(X) - S
  L <- chol(XtX.S)
  yprime <- crossprod(X,y)
  B <- backsolve(L,forwardsolve(t(L),yprime))
  B
}

N <- 5000
t1 <- rnorm(N)
t2 <- rnorm(N)
x1 <- (t1 + rnorm(N, sd = .4))
x2 <- (t2 + rnorm(N, sd = .3))
y <- 1 + 2*t1 + .5*t2 + rnorm(N)
X <- model.matrix(y ~ x1 + x2)
r <- c(1, var(t1)/var(x1), var(t2)/var(x2))
eivreg(y,X,r)

### R Code for least squares using decomposition
X <- matrix(c(1,1,1,1,10,3,15,4), ncol=2)
y <- c(5,10,15,20)
yprime <- crossprod(X,y)
L <- chol(crossprod(X))
### Step 1 solve for Z
Z <- forwardsolve(t(L),yprime)
### Step 2 solve for B. These are the least squares estimates
(B <- backsolve(L,Z))
lm(y~X[,2]) # Built in function for a check

### R Code for weighted least squares using decomposition and
elementwise calculation
X <- matrix(c(1,1,1,1,10,3,15,4), ncol=2)
y <- c(5,10,15,20)
V <- diag(runif(4)) #simulate weights on diagonal
X2 <- 1/diag(V) * X
yprime <- crossprod(X2,y)
L <- chol(crossprod(X,X2))
### Step 1 solve for Z
Z <- forwardsolve(t(L),yprime)
### Step 2 solve for B. These are the WLS estimates
```

```
(B <- backsolve(L,Z))
```

Code Examples For EAP Examples

```
library(mvtnorm)
library(statmod)

grm <- function(theta, d, score, a, D = D) {
  maxD <- length(d)
  if(score == 0) {
    pr <- 1/(1 + exp(D*a*(theta - d[(score+1)])))
  }
  else if(score == maxD) {
    pr <- 1/(1 + exp(-D*a*(theta - d[score])))
  } else {
    pr <- 1/(1 + exp(D*a*(theta - d[(score+1)])))
    - 1/(1 + exp(D*a*(theta - d[score])))
  }
  pr
}

paramList.mirt <- function(dat, dim.order = NULL){
  tmp <- split(dat, dat$strand)
  tmp <- tmp[dim.order]
  nms <- names(tmp)
  num.dimensions <- length(tmp)
  mat <- vector("list", num.dimensions)
  for(i in 1:num.dimensions){
    vars <- c('a', paste('b', 1:5, sep=''))
    poly <- subset(tmp[[i]], model == 'GRL')
    poly <- poly[, vars]
    a_vector <- as.numeric(poly$a)
    poly <- as.data.frame(t(poly[, vars[-1]]))
    py <- as.list(poly)
    py <- lapply(py, function(x) x[!is.na(x)])

    ### MC Items
    mc <- subset(tmp[[i]], model == '3pl')
    mat[[i]] <- list("3pl" = list(a = mc$a , b = mc$b, c = mc$c),
      gpcm = list(a = a_vector, d = py))
  }
  names(mat) <- dim.order
  mat
}

gauss.proc <- function(Q, num.dimensions = 4, mu, sigma, method = c(
  'gh', 'ghc', 'es', 'ghr')){
```

```

sigma <- as.matrix(sigma)
if (method == 'gh'){
  idx <- as.matrix(expand.grid(rep(list(1:Q), num.dimensions)))
  gh <- gauss.quad(Q, kind='hermite')
  nodes <- gh$nodes * sqrt(2)
  whts <- gh$weights * pi^num.dimensions/2
  nodes <- matrix(nodes[as.vector(idx)], ncol = num.dimensions)
  whts <- matrix(whts[as.vector(idx)], ncol = num.dimensions)
  whts <- whts_orig <- exp(rowSums(log(whts)))
  whts <- (dmvnorm(nodes, mean = mu, sigma = sigma)/dmvnorm(nodes,
    mean = mu, sigma = diag(1, num.dimensions))) * whts
  rotate <- FALSE
} else if (method == 'ghc'){
  idx <- as.matrix(expand.grid(rep(list(1:Q), num.dimensions)))
  gh <- gauss.quad(Q, kind='hermite')
  nodes <- gh$nodes * sqrt(2)
  whts <- gh$weights * pi^num.dimensions/2
  nodes <- matrix(nodes[as.vector(idx)], ncol = num.dimensions)
  %% chol(sigma)
  whts <- matrix(whts[as.vector(idx)], ncol = num.dimensions)
  whts <- whts_orig <- exp(rowSums(log(whts)))
  rotate <- TRUE
} else if (method == 'ghr') {
  nodes <- rmvnorm(Q, mean = mu, sigma = sigma)
  whts <- whts_orig <- 1
  idx <- NULL
  rotate <- TRUE
} else {
  idx <- as.matrix(expand.grid(rep(list(1:Q), num.dimensions)))
  nodes <- seq(-4, 4, length.out = Q)
  nodes <- matrix(nodes[as.vector(idx)], ncol = num.dimensions)
  whts <- whts_orig <- dmvnorm(nodes, mean = mu, sigma)
  rotate <- FALSE
}
list(nodes = nodes, whts = whts, idx = idx, Q = Q, rotate = rotate
, whts_orig=whts_orig)
}

mirt.eap <- function(x, params, mu, sigma, D = 1.7, dim.order = NULL
, skip = FALSE, grid, refine.steps = 1){

Q <- grid$Q
rotate <- grid$rotate
num.dimensions <- length(mu)
sigma <- as.matrix(sigma)

x <- split(x, x$strand)

```

```

scoreMat <- vector("list", num.dimensions)
names(scoreMat) <- names(x)[1:num.dimensions]
for(i in 1:length(x)){
  scoreMat[[i]] <- x[[i]][, c('score', 'model')]
}

### Must reorder so things are in same order as covariance matrix
scoreMat <- scoreMat[dim.order]

### Do error checks here
if(skip == FALSE){
  if(length(mu) != ncol(sigma)) stop('The_mean_vector_and_
    covariance_matrix_are_not_conformable')
  if(any(names(rr) != names(scoreMat))) stop('Your_scoreMat_is_
    ordered_as_', names(scoreMat),
    'but_your_item_parameters_are_ordered_as_', names(params))
  if(Q > 50) stop('You_are_using_a_LOT_of_quadrature_points._If_
    you_are_sure_you_want_to_do_this,_use_skip_=TRUE_to_bypass_
    this_message')
}

nodes <- grid$nodes
whts <- grid$whts

for(l in 1:refine.steps){ ## loop for centering

  if(l > 1) {
    SE <- sqrt(diag(vcov))
    nodes <- sweep(sweep(nodes, 2, SE, FUN='*'), 2, result, FUN='+')
    whts <- (dmvnorm(nodes, mean = mu, sigma = sigma)/dmvnorm(nodes,
      mean = rep(0,4), sigma = diag(.8,4))) * grid$whts_orig
  }

  mat <- vector("list", num.dimensions)
  for(k in 1:num.dimensions){
    a <- params[[k]]$'3pl'$a
    b <- params[[k]]$'3pl'$b
    c <- params[[k]]$'3pl'$c
    aa <- params[[k]]$gpcm$a
    d1 <- params[[k]]$gpcm$d
    x1 <- scoreMat[[k]]$score[which(scoreMat[[k]]$model == '3pl')]
    x2 <- scoreMat[[k]]$score[which(scoreMat[[k]]$model != '3pl')]
    if(rotate) {
      gh <- nodes[,k]
      Q <- length(gh)
    } else {
      gh <- unique(nodes[,k])
    }
  }
}

```

```

if(length(x1) > 0){
  tmp <- sapply(1:Q, function(i) c + (1 - c) / (1 + exp(-D * a *
    (gh[i] - b))))
  res1 <- sapply(1:Q, function(i) prod(dbinom(x1, 1, tmp[,i])))
  if(rotate){
    t1 <- res1
  } else {
    t1 <- res1[as.vector(grid$idx[,k])]
  }
} else t1 <- NULL
if(length(x2) > 0){
  res2 <- sapply(1:Q, function(i) prod((mapply(grm, d1, aa,
    theta = gh[i], score = x2, D = D))))
  if(rotate){
    t2 <- res2
  } else {
    t2 <- res2[as.vector(grid$idx[,k])]
  }
} else t2 <- NULL
mat[[k]] <- cbind(t1, t2)
}

### rn is for numerator and is likelihood * prior
rn <- apply(cbind(do.call(cbind, mat), whts), 1, prod)
### rd is 'result denominator' and is normalizing constant
rd <- sum(rn)
result <- colSums(nodes * rn)/rd
## full covariance matrix
vcov <- sapply(1:num.dimensions, function(j) sapply(1:num.
  dimensions, function(i)
    sum((nodes[,i] - result[i]) * (nodes[,j] - result[j]) * rn)/rd))

} ### end centering loop
structure(list(coefficients = result, vcov = vcov, Dimensions =
  dim.order), class = 'mirt.eap')
}

```

Code to Compute Multivariate as Product of Univariate Densities

```

library(mvtnorm)
sigma <- matrix(c(1,.5,.5,1),2,2)
L <- chol(sigma)
dmvnorm(c(1,-1), mean = c(0,0), sigma = sigma)
prod(1/diag(L)) * prod(dnorm(forwardsolve(t(L), c(1, -1))))

```

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