Spike-and-Slab LASSO Generalized Additive Models and Scalable Algorithms for High-Dimensional Data Analysis

Supporting Information

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Supplementary Information 1: Marginal Distribution of γ_i^*

Given that $\gamma_j^*|\gamma_j, \theta_j \sim Bin(1, \gamma_j\theta_j)$ where $\gamma_j|\theta_j \sim Bin(1, \theta_j)$, we can derive the marginal distribution of γ_j^* with the following manipulation.

$$\begin{split} &\Pr(\gamma_j^*=1|\theta_j) = \Pr(\gamma_j^*=1,\gamma_j=1|\theta_j) + \Pr(\gamma_j^*=1,\gamma_j=0|\theta_j) \\ &= \Pr(\gamma_j^*=1,\gamma_j=1|\theta_j) + 0 \quad \text{[hierarchical structure between } \gamma^* \text{ and } \gamma.] \\ &= \Pr(\gamma_j^*=1|\gamma_j=1,\theta_j) \Pr(\gamma_j=1|\theta_j) \\ &= \theta_j^2 \\ &\Pr(\gamma_j^*=0|\theta_j) = \Pr(\gamma_j^*=0,\gamma_j=1|\theta_j) + \Pr(\gamma_j^*=0,\gamma_j=0|\theta_j) \\ &= \Pr(\gamma_j^*=0,\gamma_j=1|\theta_j) + \Pr(\gamma_j^*=0,\gamma_j=0|\theta_j) \\ &= \Pr(\gamma_j^*=0|\gamma_j=1,\theta_j) \Pr(\gamma_j=1|\theta_j) + \Pr(\gamma_j^*=0|\gamma_j=0,\theta_j) \Pr(\gamma_j=0|\theta_j) \\ &= (1-\theta_j)\theta_j + 1(1-\theta_j) = 1-\theta_j^2 \end{split}$$

Supplementary Information 2: EM-Iterative Weighted Least Square Algorithm for BHAM Model Fitting

Similar to the EM-CD algorithm, the EM-Iterative Weighted Least Square (EM-IWLS) algorithm is an EM-based algorithm where the iterative weighted least squares algorithm is used to find the estimate of β , ϕ that maximizes $E(Q_1)$. The iterative weighted least squares algorithm was originally proposed to fit the classical generalized linear models, and generalized to fit some Bayesian hierarchical models. (Gelman et al. 2013) Yi and Ma (Yi and Ma 2012) extended the algorithm to fit Bayesian hierarchical models for high-dimensional data analysis. Specifically, the authors formulated Student's t-distribution and double exponential distribution as hierarchical normal distributions such that generalized linear models with shrinkage priors can be easily fitted. In this work, we further extend the EM-IWLS paradigm to fit the proposed BHAM method. Compare to the EM-CD algorithm, EM-IWLS estimates the variance-covariance matrix of the coefficients, providing an opportunity to derive the uncertainty quantification of smooth functions. We will defer this discussion to future work, due to the delicacy of the topic. However, the EM-IWLS is not as computationally efficient as EM-CD, particularly in high-dimensional settings.

A double exponential prior, $\beta | S \sim DE(0, S)$ can be formulated as a hierarchical normal prior with unknown variance τ^2 integrated out:

$$\beta | \tau^2 \sim N(0, \tau^2)$$

 $\tau^2 | S \sim Gamma(1, 1/(2S^2)),$

For the mixture double exponential priors, we can define the scale parameter $S = (1 - \gamma)s_0 + \gamma s_1$. The change in the prior formulation in turn leads to the change in the log posterior density function, as Q_1 needs to account for the hyperprior of τ^2 :

$$Q_1(\boldsymbol{\beta}, \phi) = \log f(\mathbf{y}|\boldsymbol{\beta}, \phi) + \sum_{j=1}^{p} \left[\log f(\beta_j | \tau_j^2) + \log f(\tau_j^2 | S_j) + \sum_{k=1}^{K_j} \{ \log f(\beta_{jk}^* | \tau_{jk}^{*2}) + \log f(\tau_{jk}^{*2} | S_j^*) \} \right]. \quad (1)$$

Since τ^2 are not of our primary interest, we treat them as the "missing" data in addition to the latent indicators γ , and hence construct the expectation $E_{\gamma,\tau^2|\Theta^{(t-1)}}(Q_1)$ in the E-step. To note, unlike the same latent indicator γ_j^* which is shared by the coefficients of the nonlinear terms β_{jk}^* for $k=1,\ldots,K_j$, τ_{jk}^2 is coefficient specific for β_{jk}^* . $E(S_j^{-1}|\beta_j,s_0,s_1)$, $E(S_j^{*-1}|\beta_j^*,s_0,s_1)$, $E(\tau_j^2|S_j,\beta_j)$ and $E(\tau_{jk}^{*2}|S_j^*,\beta_{jk}^*)$ needs to be calculated to formulate $E(Q_1)$. As neither $E(S_j^{-1}|\beta_j,s_0,s_1)$ nor $E(S_j^{*-1}|\beta_j^*,s_0,s_1)$ depends on τ^2 s, they can be derived following the same derivation in the EM-CD algorithm. On the other hand, τ^2 , following gamma distributions, is a conjugate prior for the normal variance, and the conditional posterior density of τ^{-2} is an inverse Gaussian distribution. $E(\tau_j^{-2})$ and $E(\tau_{jk}^{*-2})$ are calculated using the closed form equation

$$E(\tau_j^{-2}|S_j,\beta_j) = S_j^{-1}/|\beta_j|$$
 $E(\tau_{jk}^{*-2}|S_j^*,\beta_{jk}^*) = S_j^{*-1}/|\beta_{jk}^*|,$

where S_j and S_j^* are replaced by the expectation and β s are replaced with $\beta^{(t-1)}$. With simplification (up to constant additive terms), we have

$$E(Q_1) = \log f(\mathbf{y}|\boldsymbol{\beta}, \phi) - \sum_{j=1}^{p} \left[2E(\tau_j^{-2})\beta_j^2 + \sum_{k=1}^{K_j} 2E(\tau_{jk}^{*-2})\beta_{jk}^{*-2} \right].$$
 (2)

 $2E(\tau^{-2})\beta^2$ can be seen as the kernel of a normal density with mean 0 and variance $E(\tau^2)$, and we can formulate the coefficients β as a multivariate normal distribution with means **0** and variance covariance matrix Σ_{τ^2} , where Σ_{τ^2} is a diagonal matrix with $E(\tau^2)$ s on the diagonal,

$$\boldsymbol{\beta} \sim \text{MVN}(0, \boldsymbol{\Sigma}_{\tau^2}).$$

Meanwhile, following the classical IWLS, we can approximate the generalized model likelihood at each iteration with a weighted normal likelihood:

$$f(\mathbf{y}|\boldsymbol{\beta}, \phi) \approx \text{MVN}(\mathbf{z}|\boldsymbol{X}\boldsymbol{\beta}, \phi\boldsymbol{\Sigma})$$

where the 'normal response' z_i and 'weight' w_i are called the pseudo-response and pseudo-weight respectively. The pseudo-response and the pseudo-weight are calculated by

$$z_i = \hat{\eta}_i - rac{L^{'}(y_i|\hat{\eta}_i)}{L^{''}(y_i|\hat{\eta}_i)} \quad w_i = -L^{''}(y_i|\hat{\eta}_i),$$

where $\hat{\eta}_i = (\mathbf{X}\hat{\boldsymbol{\beta}})_i$, $L'(y_i|\hat{\eta}_i,\hat{\phi})$ and $L''(y_i|\hat{\eta}_i,\hat{\phi})$ are the first and second derivative of the log density, $\log f(\mathbf{y}_i|\boldsymbol{\beta},\phi)$ with respect to η_i .

With $z \sim \text{MVN}(X\beta, \phi \Sigma)$ and $\beta \sim \text{MVN}(0, \phi \Sigma_{\tau^2})$, we can augment the two multivariate normal distributions and update the estimates for β and ϕ via least squares in each iteration of the EM algorithm. We create the augmented response, augmented data, and augmented variance-covariance matrix following

$$oldsymbol{z}_* = egin{bmatrix} oldsymbol{z} \ 0 \end{bmatrix} \hspace{1cm} oldsymbol{X}_* = egin{bmatrix} oldsymbol{X} \ I \end{bmatrix} \hspace{1cm} oldsymbol{\Sigma}_* = egin{bmatrix} oldsymbol{\Sigma} \ 0 & oldsymbol{\Sigma}_{ au^2}/\phi \end{bmatrix},$$

such that

$$z_* \sim \text{MVN}(X_*\beta, \phi\Sigma_*).$$

Using the least squares estimators to update β and ϕ , we have

$$\boldsymbol{\beta}^{(t)} = (\boldsymbol{X}_*^T \boldsymbol{\Sigma}^{-1} \boldsymbol{X}_*)^{-1} \boldsymbol{X}_*^T \boldsymbol{\Sigma}^{-1} \boldsymbol{z}_* \qquad \qquad \boldsymbol{\phi}^{(t)} = \frac{1}{n} (\boldsymbol{z}_* - X_* \boldsymbol{\beta}^{(t)})^T \boldsymbol{\Sigma}^{-1} (\boldsymbol{z}_* - X_* \boldsymbol{\beta}^{(t)}).$$

To note, the variance-covariance matrix of the coefficient estimates variance-covariance matrix can be derived in the EM-IWLS algorithm and in turn can be used for statistical inferences,

$$\operatorname{Var}(\boldsymbol{\beta}^{(t)}) = (\boldsymbol{X}_*^T \boldsymbol{\Sigma}^{-1} \boldsymbol{X}_*)^{-1} \phi^{(t)}.$$

Totally, the proposed EM-IWLS algorithm is summarized as follows:

- 1) Choose a starting value $\boldsymbol{\beta}^{(0)}$ and $\boldsymbol{\theta}^{(0)}$ for $\boldsymbol{\beta}$ and $\boldsymbol{\theta}$. For example, we can initialize $\boldsymbol{\beta}^{(0)} = \mathbf{0}$ and $\boldsymbol{\theta}^{(0)} = \mathbf{0}$.
- 2) Iterate over the E-step and M-step until convergence

E-step: calculate $E(\gamma_j)$, $E(\gamma_j^*)$ and $E(\tau_j^{-2})$, $E(\tau_{jk}^{*-2})$ with the estimates $\Theta^{(t-1)}$ from the previous iteration

M-step:

- a) Based on the current value of β , calculate the pseudo-data $z_i^{(t)}$ and the pseudo-weights $w_i^{(t)}$
- b) Update $\beta^{(t)}$ by running the augmented weighted least squared
- c) If ϕ is present, update ϕ

Similar to EM-CD, we assess convergence by the criterion, $|d^{(t)} - d^{(t-1)}|/(0.1 + |d^{(t)}|) < \epsilon$, where ϵ is a small value (say 10^{-5}).

Supplementary Information 3: Generalization of BHAM framework

In the manuscript, we describe the Bayesian hierarchical additive model with the two-part spike-and-slab LASSO prior. Nevertheless, the proposed model and algorithm can be easily generalized to accommodate other priors thanks to the reparameterization of smooth functions. To be more specific, we can apply a regularized prior on the linear coefficient, and a group regularized prior for the nonlinear coefficients. For example, we can apply the same proposed framework with a spike-and-slab mixture normal prior,

$$\beta_j | \gamma_j, s_0, s_1 \sim N(0, (1 - \gamma_j) s_0 + \gamma_j s_1)$$

$$\beta_{jk}^* | \gamma_j^*, s_0, s_1 \stackrel{\text{iid}}{\sim} N(0, (1 - \gamma_j^*) s_0 + \gamma_j^* s_1), k = 1, \dots, K_j.$$

The algorithm derivation still follows with slight modification, replacing l_1 penalization with l_2 penalization. The implementation of effect hierarchy would be more challenging for the priors that do rely on a latent indicator for variable selection. As a naive solution, readers can consider the linear prior and the nonlinear prior being independent at the cost of bi-level selection accuracy.

Supplementary Information 4: Predictive Performance of Linear Simulations

P	mgcv	LASSO	COSSO	Adaptive COSSO	BHAM	SB-GAM	spikeSlabGAM
4	0.38 (0.01)	0.39 (0.01)	0.31 (0.08)	0.29 (0.11)	0.38 (0.01)	0.35(0.01)	0.39 (0.01)
10	0.36(0.02)	0.38(0.01)	0.35(0.03)	0.34 (0.04)	0.39(0.01)	0.33(0.02)	0.39(0.01)
50	0.09(0.09)	0.37(0.01)	$0.30 \ (0.06)$	$0.30 \ (0.36)$	0.38(0.01)	0.32(0.03)	0.37(0.01)
100	_	0.37(0.01)	0.28(0.07)	0.34 (0.04)	0.38(0.01)	0.29(0.07)	0.35(0.01)
200	_	0.36(0.01)	0.26(0.08)	0.31(0.06)	0.38(0.03)	0.28(0.06)	0.33(0.02)

Table 1: The average and standard deviation of the out-of-sample R^2 measure for Gaussian outcomes over 50 iterations. The models of comparison include the proposed Bayesian hierarchical additive model (BHAM), linear LASSO model (LASSO), component selection and smoothing operator (COSSO), adaptive COSSO, mgcv, sparse Bayesian generalized additive model (SB-GAM), and spikeSlabGAM model. mgcv doesn't provide estimation when number of parameters exceeds sample size i.e. p = 100, 200.

P	mgcv	LASSO	COSSO	Adaptive COSSO	BHAM	SB-GAM	spikeSlabGAM
4	0.79 (0.01)	0.79 (0.01)	0.76 (0.04)	0.75 (0.04)	0.78 (0.01)	0.76(0.01)	0.79 (0.01)
10	0.77(0.01)	0.79(0.01)	0.78(0.01)	0.78(0.01)	0.78(0.01)	0.75(0.01)	0.79(0.01)
50	0.62(0.01)	0.78(0.01)	0.75(0.03)	0.73(0.04)	0.74(0.07)	0.75(0.02)	0.77(0.01)
100	_	0.78(0.01)	0.73(0.04)	0.69 (0.05)	0.73(0.07)	0.74(0.02)	0.76(0.02)
200	-	0.78(0.01)	0.71(0.05)	0.67(0.05)	0.73(0.06)	0.73(0.03)	0.72(0.03)

Table 2: The average and standard deviation of the out-of-sample area under the curve measures for binomial outcomes over 50 iterations. The models of comparison include the proposed Bayesian hierarchical additive model (BHAM), linear LASSO model (LASSO), component selection and smoothing operator (COSSO), adaptive COSSO, mgcv, sparse Bayesian generalized additive model (SB-GAM), and spikeSlabGAM model. mgcv doesn't provide estimation when number of parameters exceeds sample size i.e. p = 100, 200.

Supplementary Information 5: Variable Selection Performance of Simulations

P	Metric	LASSO	COSSO	Adaptive COSSO	BHAM	SB-GAM	spikeSlabGAM
4	Precision	1.00	1.00	1.00	1.00	1.00	1.00
10	Precision	0.58	0.71	0.69	0.93	0.88	0.89
50	Precision	0.38	0.60	0.59	0.77	0.80	0.52
100	Precision	0.35	0.62	0.65	0.82	0.77	0.42
200	Precision	0.29	0.65	0.57	0.21	0.74	0.36
4	Recall	0.54	0.61	0.54	0.38	0.97	0.55
10	Recall	0.46	0.50	0.54	0.38	0.97	0.54
50	Recall	0.34	0.30	0.30	0.32	0.90	0.55
100	Recall	0.27	0.30	0.25	0.27	0.92	0.54
200	Recall	0.27	0.25	0.29	0.56	0.92	0.53
10	MCC	0.18	0.35	0.36	0.47	0.86	0.55
50	MCC	0.25	0.34	0.36	0.45	0.83	0.47
100	MCC	0.24	0.37	0.37	0.45	0.82	0.43
200	MCC	0.23	0.36	0.36	0.23	0.81	0.40

Table 3: The variable selection performance of binomial simulations, measured by positive predictive value (precision), true positive rate (recall), and Matthews correlation coefficient (MCC), for the high-dimensional methods averaged over 50 iterations. The models of comparison include the proposed Bayesian hierarchical additive model (BHAM), linear LASSO model (LASSO), component selection and smoothing operator (COSSO), adaptive COSSO, sparse Bayesian generalized additive model (SB-GAM), and spikeSlabGAM model. MCC is ill-defined when p=4 simulation (no true negative), and hence omitted for all methods.

P	Metric	LASSO	COSSO	Adaptive COSSO	BHAM	SB-GAM	spikeSlabGAM
4	Precision	1.00	1.00	1.00	1.00	1.00	1.00
10	Precision	0.59	0.97	0.97	0.90	0.56	0.99
50	Precision	0.43	0.74	0.84	0.89	0.39	0.99
100	Precision	0.34	0.53	0.72	0.91	0.29	0.99
200	Precision	0.27	0.42	0.52	0.92	0.33	0.99
4	Recall	1.00	0.87	0.84	1.00	1.00	1.00
10	Recall	1.00	0.98	0.96	1.00	1.00	1.00
50	Recall	1.00	0.85	0.99	1.00	1.00	1.00
100	Recall	1.00	0.84	0.97	1.00	1.00	1.00
200	Recall	1.00	0.78	0.88	1.00	0.98	1.00
10	MCC	0.59	0.96	0.94	0.91	0.54	0.99
50	MCC	0.59	0.76	0.90	0.92	0.57	1.00
100	MCC	0.54	0.61	0.81	0.94	0.48	1.00
200	MCC	0.49	0.52	0.64	0.95	0.52	0.99

Table 4: The variable selection performance of linear Gaussian simulations, measured by positive predictive value (precision), true positive rate (recall), and Matthews correlation coefficient (MCC), for the high-dimensional methods averaged over 50 iterations. The models of comparison include the proposed Bayesian hierarchical additive model (BHAM), linear LASSO model (LASSO), component selection and smoothing operator (COSSO), adaptive COSSO, sparse Bayesian generalized additive model (SB-GAM), and spikeSlabGAM model. MCC is ill-defined when p=4 simulation (no true negative), and hence omitted for all methods.

P	Metric	LASSO	COSSO	Adaptive COSSO	BHAM	SB-GAM	spikeSlabGAM
4	Precision	1.00	1.00	1.00	1.00	1.00	1.00
10	Precision	0.61	0.97	0.98	0.55	0.74	0.91
50	Precision	0.35	0.59	0.68	0.26	0.59	0.61
100	Precision	0.28	0.47	0.51	0.30	0.46	0.57
200	Precision	0.26	0.44	0.42	0.38	0.41	0.38
4	Recall	1.00	0.85	0.78	1.00	1.00	1.00
10	Recall	1.00	0.96	0.97	0.99	1.00	1.00
50	Recall	1.00	0.88	0.74	0.92	1.00	1.00
100	Recall	1.00	0.78	0.56	0.89	1.00	0.99
200	Recall	1.00	0.67	0.48	0.89	0.99	0.98
10	MCC	0.58	0.95	0.96	0.49	0.72	0.91
50	MCC	0.52	0.66	0.65	0.37	0.73	0.74
100	MCC	0.49	0.55	0.47	0.43	0.65	0.72
200	MCC	0.49	0.49	0.39	0.47	0.62	0.58

Table 5: The variable selection performance of linear binomial simulations, measured by positive predictive value (precision), true positive rate (recall), and Matthews correlation coefficient (MCC), for the high-dimensional methods averaged over 50 iterations. The models of comparison include the proposed Bayesian hierarchical additive model (BHAM), linear LASSO model (LASSO), component selection and smoothing operator (COSSO), adaptive COSSO, sparse Bayesian generalized additive model (SB-GAM), and spikeSlabGAM model. MCC is ill-defined when p=4 simulation (no true negative), and hence omitted for all methods.

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