CPSC 540: Assignment 2

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1 Convergence Rates

1.1 Gradient Descent

- 1. If f is a C^1 function, strong convexity means $f(y) \geq f(x) + \langle f'(y), y x \rangle + \frac{u}{2}||x y||^2$ for u > 0 and y, x in its domain. Applying this inequality at x^t and x^* , we have $f(x^t) \geq f(x^*) + \langle f'(x^*), x^* x^T \rangle + \frac{u}{2}||x^* x^t||^2$. Since $f'(x^*) = 0$, $\frac{u}{2}||x^* x^t||^2 \leq f(x^t) f(x^*) = O(\rho^t)$. This means the iteration $||x^t x^*||$ has a convergence rate of $O(\rho^{t/2})$. If f is a C^2 function, by Taylor's theorem, there is a z such that $f(x^t) = f(x^*) + \nabla f(x^*)(x^t x^*) + \frac{1}{2}(x^t x^*)^T \nabla^2 f(u)(x^t x^*)$. By strong convexity, $f(x^*) \geq f(x^t) + \nabla f(x^*)(x^t x^*) + \frac{u}{2}||x^t x^*||^2$. Again, as $\nabla f(x^*) = 0$, we have $\frac{u}{2}||x^* x^t||^2 \leq f(x^t) f(x^*) = O(\rho^t)$, which means the iteration has a convergence rate of $O(\rho^{t/2})$.
- 2. From the descent lemma, we know $f(x^{t+1}) \leq f(x^t) + \nabla f(x^t)^T (x^{t+1} x^t) + \frac{L}{2} ||x^{t+1} x^t||^2$. With the constant step size, $x^{t+1} x^t = \alpha \nabla f(x^t)$. Plugging this into the descent lemma, we get

$$f(x^{t+1}) \le f(x^t) - \alpha \nabla f(x^t)^T \nabla f(x^t) + \frac{L\alpha^2}{2} ||\nabla f(x^t)||^2$$
$$= f(x^t) - (\alpha - \frac{L\alpha^2}{2}) ||\nabla f(x^t)||^2$$

Comparing the upper bound with the proof in class, we define $\frac{1}{2\tilde{L}} = \alpha - \frac{L\alpha^2}{2}$ and only need to show $\tilde{L} \geq L$ in order to preserve the linear convergence rate. Since $\alpha < \frac{L}{2}$, the minimum value of \tilde{L} is L. Thus, we show that the convergence rate is linear when using a constant step size.

3. By descent lemma, we have

$$\begin{split} f(x^{t+1}) &\leq f(x^t) - \frac{1}{L^t} \nabla f(x^t)^T \nabla f(x^t) + \frac{L}{2L^{t2}} ||\nabla f(x^t)||^2 \\ &= f(x^t) - (\frac{1}{L^t} - \frac{L}{2L^{t2}}) ||\nabla f(x^t)||^2 \end{split}$$

To satisfy the inequality, we need $\frac{1}{L^t} - \frac{L}{2L^{t2}} \ge \frac{1}{2L^t}$ which gives us $L^t \ge L$. Since we start with some L^0 that is smaller than L and double the value when the inequality is not

satisfied, $L^t \leq 2L$ always holds. Following the steps in class, we have $f(x^t) - f(x^*) \leq (1 - \frac{u}{L^0})(1 - \frac{u}{L^1}) \cdots (1 - \frac{1}{L^{t-1}})(f(x^0) - f(x^*)) \leq (1 - \frac{u}{2L})^t (f(x^0) - f(x^*))$.

4. When $L^t > L$ for any t, the convergence rate would be faster than $\rho = 1 - \frac{u}{L}$.

1.2 Sign-Based Gradient Descent

1. By the Lipschitz continuity in ∞ -norm, we have

$$\begin{split} f(x^{t+1}) &\leq f(x^t) + \nabla f(x^t)^T (x^{t+1} - x^t) + \frac{L_{\infty}}{2} \|x^{t+1} - x^t\|_{\infty}^2 \\ &= f(x^t) + \nabla f(x^t)^T \Big(-\frac{\|\nabla f(x^t)\|_1}{L_{\infty}} sign(\nabla f(x^t)) \Big) + \frac{L_{\infty}}{2} \Big\| -\frac{\|\nabla f(x^t)\|_1}{L_{\infty}} sign(\nabla f(x^t)) \Big\|_{\infty}^2 \\ &= f(x^t) - \frac{1}{L_{\infty}} \|\nabla f(x^t)\|_1^2 + \frac{1}{2L_{\infty}} \|\nabla f(x^t)\|_1^2 \\ &= f(x^t) - \frac{1}{2L_{\infty}} \|\nabla f(x^t)\|_1^2 \\ &\leq f(x^t) - \frac{1}{2L_{\infty}} \|\nabla f(x^t)\|_2^2 \end{split}$$

By the strong convexity, we have

$$-||\nabla f(x^t)||_2^2 \le -2\mu(f(x^t) - f(x^*)).$$

Therefore, we get $f(x^{t+1}) - f(x^*) \le (1 - \frac{u}{L_{\infty}})(f(x^t) - f(x^*)).$

2. If f is L_{∞} Lipschitz continuous,

$$||f'(x) - f'(y)||_2 \le ||f'(x) - f'(y)||_1$$

 $\le L_{\infty} ||x - y||_{\infty}$
 $\le L_{\infty} ||x - y||_2$

As L is the minimum value of the Lipschitz L_2 -norm constant, we have $L \leq L_{\infty}$. Similarly, if f is L_2 Lipschitz continuous, we have

$$||f'(x) - f'(y)||_{1} \leq \sqrt{d}||f'(x) - f'(y)||_{2}$$

$$\leq \sqrt{d}L||x - y||_{2}$$

$$\leq \sqrt{d}L\sqrt{d}||x - y||_{\infty}$$

$$\leq dL||x - y||_{\infty}$$

As L_{∞} is the minimum value of the Lipschitz L_{∞} -norm constant, we have $L_{\infty} \leq dL$. In conclusion, $L \leq L_{\infty} \leq dL$.

1.3 Block Coordinate Descent

1. By block-wise strong-smoothness and $p(b_t = b) = \frac{1}{k}$, we have

$$f(x^{t+1}) \leq f(x^{t}) - \frac{1}{L} \nabla_{b_{t}} f(x^{t})^{T} (\nabla f(x^{t}) \circ e_{b_{t}}) + \frac{1}{2L} ||\nabla f(x^{t}) \circ e_{b_{t}}||^{2}$$

$$= f(x^{t}) - \frac{1}{2L} ||\nabla f(x^{t}) \circ e_{b_{t}}||^{2}$$

$$\mathbf{E}[f(x^{t+1})] \leq \mathbf{E}[f(x^{t}) - \frac{1}{2L} ||\nabla f(x^{t}) \circ e_{b_{t}}||^{2}]$$

$$= \frac{1}{k} \sum_{b=1}^{k} (f(x^{t}) - \frac{1}{2L} ||\nabla f(x^{t}) \circ e_{b_{t}}||^{2})$$

$$= f(x^{t}) - \frac{1}{2kL} \sum_{b=1}^{k} ||\nabla f(x^{t}) \circ e_{b_{t}}||^{2}$$

$$= f(x^{t}) - \frac{1}{2kL} ||\nabla f(x^{t})||^{2}$$

From strong convexity, we can re-write the inequality into:

$$\mathbf{E}(f(x^{t+1})) - f(x^*) \le f(x^t) - f(x^*) - \frac{u}{kL} [f(x^t) - f(x^*)]$$
$$= (1 - \frac{u}{kL}) [f(x^t) - f(x^*)]$$

2. Assuming that each block b has its own strong-smoothness constant L_b and the blocks are sampled proportional to L_{b_t} , we have

$$f(x^{t+1}) \leq f(x^{t}) - \frac{1}{L_{b_{t}}} \nabla_{b_{t}} f(x^{t})^{T} (\nabla f(x^{t}) \circ e_{b_{t}}) + \frac{1}{2L_{b_{t}}} ||\nabla f(x^{t}) \circ e_{b_{t}}||^{2}$$

$$= f(x^{t}) - \frac{1}{2L_{b_{t}}} ||\nabla f(x^{t}) \circ e_{b_{t}}||^{2}$$

$$\mathbf{E}[f(x^{t+1})] \leq \mathbf{E}[f(x^{t}) - \frac{1}{2L_{b_{t}}} ||\nabla f(x^{t}) \circ e_{b_{t}}||^{2}]$$

$$= \sum_{b=1}^{k} \frac{L_{b_{t}}}{\sum L_{b_{t}}} (f(x^{t}) - \frac{1}{2L_{b_{t}}} ||\nabla f(x^{t}) \circ e_{b_{t}}||^{2})$$

$$= f(x^{t}) - \sum_{b=1}^{k} \frac{1}{2\sum L_{b_{t}}} ||\nabla f(x^{t}) \circ e_{b_{t}}||^{2}$$

$$\leq f(x^{t}) - \frac{u}{\sum L_{b_{t}}} [f(x^{t}) - f(x^{*})]$$

$$= (1 - \frac{u}{\sum L_{b_{t}}}) [f(x^{t}) - f(x^{*})]$$

The second last line is from applying the strong convexity block-wise. Since $L = \max_{b_t} \{L_{b_t}\}$, $L_{b_t} \leq L$, $\sum L_{b_t} \leq kL$. If there exists some $L_{b_t} \neq L$, $1 - \frac{u}{\sum L_{b_t}} \leq 1 - \frac{u}{kL}$, meaning we have a faster convergence rate.

2 Large-Scale Algorithms

2.1 Coordinate Optimization

1. The final function value is 1.4724e+02 and the total time is 2.328713 seconds.

```
function [model] = logisticL2(X, y, lambda)
  % Add bias variable
  [n,d] = size(X);
X = [ones(n,1) X];
  d = d+1;
  % Initial values of regression parameters
  w = zeros(d,1);
10
  % Optimization parameters
   maxPasses = 500;
   progTol = 1e-4;
  L = .25*max(sum(X.^2, 1)) + lambda;
14
15
   w_{-}old = w;
16
17
  Xw = X*w;
18
19
   for t = 1: maxPasses*d
20
21
       % Choose variable to update 'j'
22
       j = randi(d);
23
24
       % Compute partial derivative 'g_j'
25
26
       yXw = y.*(Xw);
27
       sigmoid = 1./(1 + \exp(-yXw));
28
       g_{-j} = -(y.*(1-sigmoid))'*X(:,j) + lambda*w(j);
29
30
       % Update variable
31
       w(j) = w(j) - (1/L) * g_{-j};
32
       Xw = Xw - (1/L)*X(:, j)*g_{-j};
33
34
       % Check for lack of progress after each "pass"
35
       if \mod(t,d) == 0
36
            change = norm(w-w_old, inf);
37
            fprintf('Passes = %d, function = %.4e, change = %.4f\n',t/d,
38
               logisticL2\_loss(w,X,y,lambda),change);
            if change < progTol
39
                fprintf('Parameters changed by less than progTol on pass\n');
40
                break;
41
```

```
end
42
            w_{-}old = w;
43
       end
44
  end
45
46
47
  model.w = w;
48
   model.predict = @predict;
49
50
2. The final function value is 1.4418e+02 and the total number of passes is 153.
   function [model] = logisticL2(X,y,lambda)
  % Add bias variable
  [n,d] = size(X);
  X = [ones(n,1) X];
  d = d+1;
  % Initial values of regression parameters
8
  w = zeros(d,1);
9
10
  % Optimization parameters
   maxPasses = 500;
12
   progTol = 1e-4;
   L_candidate = .25*(sum(X.^2, 1)) + lambda;
14
   w_old = w;
16
  Xw = X*w;
17
18
   for t = 1: maxPasses*d
19
20
       % Choose variable to update 'j'
21
       \%j = randi(d);
22
23
       p = L_candidate/(sum(L_candidate));
24
       j = sampleDiscrete(p);
25
       L = L_{candidate(j)};
26
27
       % Compute partial derivative 'g_j'
29
       yXw = y.*(Xw);
30
       sigmoid = 1./(1+\exp(-yXw));
31
       g_{-j} = -(y.*(1-sigmoid))'*X(:,j) + lambda*w(j);
32
33
       %g_{-j} = g(j);
34
35
       % Update variable
36
       w(j) = w(j) - (1/L) * g_{-j};
37
       Xw = Xw - (1/L)*X(:,j)*g_{-j};
38
39
       % Check for lack of progress after each "pass"
40
       if \mod(t,d) == 0
41
            change = norm(w-w_old, inf);
42
            fprintf('Passes = %d, function = %.4e, change = %.4f\n',t/d,
43
```

```
logisticL2_loss(w,X,y,lambda),change);
            if change < progTol
44
                fprintf('Parameters changed by less than progTol on pass\n');
45
46
            end
47
            w_old = w;
48
       end
49
   end
50
51
   model.w = w:
53
   model.predict = @predict;
55
3. The final function value is 1.4091e+02 and the total number of passes is 149.
   function [model] = logisticL2(X, y, lambda)
  % Add bias variable
   [n,d] = size(X);
  X = [ones(n,1) X];
  d = d+1;
  % Initial values of regression parameters
8
  w = zeros(d,1);
9
10
  % Optimization parameters
   maxPasses = 500;
12
   progTol = 1e-4;
   L_candidate = .25*(sum(X.^2, 1)) + lambda;
   w_{-}old = w;
15
16
  Xw = X*w;
17
18
   for t = 1: maxPasses*d
19
20
       % Choose variable to update 'j'
21
       j = randi(d);
22
       L = L_{-candidate(j)};
23
       % Compute partial derivative 'g_j'
25
26
       yXw = y.*(Xw);
27
       sigmoid = 1./(1+exp(-yXw));
       g_{-j} = -(y.*(1-sigmoid)) *X(:,j) + lambda*w(j);
29
30
       % Update variable
31
       w(j) = w(j) - (1/L)*g_{-j};
32
       Xw = Xw - (1/L)*X(:, j)*g_{-j};
33
34
       % Check for lack of progress after each "pass"
35
       if mod(t,d) == 0
36
            change = norm(w-w_old, inf);
37
            fprintf('Passes = %d, function = %.4e, change = %.4f\n',t/d,
38
                logisticL2_loss(w,X,y,lambda),change);
```

```
if change < progTol
39
                 fprintf('Parameters changed by less than progTol on pass\n');
40
41
            end
42
            w_{old} = w;
43
44
       end
   end
45
46
47
   model.w = w;
   model.predict = @predict;
49
   end
```

4. We stated in class that L_{j_t} is used as the step size and sampling j_t proportional to L_{j_t} gives

$$E[f(x^t)] - f(x^t) \le (1 - \frac{\mu}{d\bar{L}})[f(x^0) - f(x^*)].$$

By using the uniform sampling, $E[\bar{L}]_{Uniform} = \frac{1}{d} \sum_{i=1}^{d} L_i$

By using the Lipschitz sampling, $E[\bar{L}]_{Lipschitz} = \sum_{i=1}^d p_i L_i = \frac{1}{\sum_{i=1}^d L_i} \sum_{i=1}^d L_i^2$

To prove that uniform sampling has a tighter bound than Lipschitz sampling, we need to show that

$$E[\bar{L}]_{Uniform} \le E[\bar{L}]_{Lipschitz},$$

which is equivalent to showing

$$\frac{1}{d} \left(\sum_{i=1}^d L_i \right)^2 \le \sum_{i=1}^d L_i^2.$$

According to the CauchySchwarz inequality, the above inequality holds and thus uniform sampling outperforms Lipschitz sampling.

2.2 Proximal-Gradient

1. errTrain = 0.0060; errTest = 0.2740 The number of non-zero parameters: 500

The number of original features that the model uses: 100

```
function [model] = softmaxClassifierL2(X,y,lambda)

% Compute sizes

[n,d] = size(X);

k = max(y);

W = zeros(d,k); % Each column is a classifier

W(:) = findMin(@softmaxLoss,W(:),500,1,X,y,k,lambda);

model.W = W;

model.predict = @predict;
```

```
end
13
   function [yhat] = predict (model, X)
  W = model.W;
   [ , yhat ] = max(X*W, [], 2);
16
17
18
   function [nll,g,H] = softmaxLoss(w,X,y,k,lambda)
19
20
   [n,p] = size(X);
^{21}
  W = reshape(w, [p k]);
22
23
  XW = X*W;
24
   Z = sum(exp(XW), 2);
25
26
   ind = sub2ind([n k], [1:n]', y);
   n11 = -sum(XW(ind) - log(Z)) + sum(sum((lambda/2)*W.^2));
28
30
   g = zeros(p,k);
31
   for c = 1:k
32
       g(:,c) = X'*(exp(XW(:,c))./Z-(y == c)) + lambda*W(:,c);
33
   end
34
35
   g = reshape(g, [p*k 1]);
2. The number of non-zero parameters: 35
   The number of original features that the model uses: 19
   function [model] = softmaxClassifierL1(X,y,lambda,maxIter)
  % Compute sizes
   [n,d] = size(X);
   k = \max(y);
  W = zeros(d,k); % Each column is a classifier
  W(:) = proxGradL1(@softmaxLoss,W(:),X,y,k,lambda,maxIter);
10
   model.W = W;
11
   model.predict = @predict;
13
14
   function [yhat] = predict (model, X)
15
  W = model.W;
   [ , yhat ] = max(X*W, [], 2);
17
19
20
   function [nll,g] = softmaxLoss(w,X,y,k,lambda)
21
   [n,p] = size(X);
^{23}
  W = reshape(w, [p k]);
24
25
```

```
26 \text{ XW} = \text{X*W};
   Z = sum(exp(XW), 2);
27
28
   ind = sub2ind([n k],[1:n]',y);
29
   nll = -sum(XW(ind) - log(Z));
30
31
32
   g = zeros(p,k);
   for c = 1:k
33
       g(:,c) = X'*(exp(XW(:,c))./Z-(y == c));
34
35
   g = reshape(g, [p*k 1]);
36
37
   end
3. errTrain = 0.0220; errTest = 0.0540 The number of non-zero parameters: 115
   The number of original features that the model uses: 23
   function [model] = softmaxClassiferGL1(X, y, lambda, maxIter)
  % Compute sizes
   [n,d] = size(X);
  k = \max(y);
  W = zeros(d,k); % Each column is a classifier
8
  W(:) = proxGradGroupL1 (@softmaxLoss,W(:),X,y,k,lambda,maxIter);
10
   model.W = W;
11
   model.predict = @predict;
12
   end
13
14
   function [yhat] = predict (model, X)
  W = model.W;
16
   [ , yhat ] = max(X*W, [], 2);
17
   end
18
19
20
   function [nll,g] = softmaxLoss(w,X,y,k,lambda)
21
22
   [n,p] = size(X);
23
  W = reshape(w, [p k]);
^{24}
25
  XW = X*W;
26
   Z = sum(exp(XW), 2);
27
28
   ind = sub2ind([n k],[1:n]',y);
29
30
   tmp = sqrt(sum(W.^2, 2));
31
32
   nll = -sum(XW(ind) - log(Z));
33
34
   g = zeros(p,k);
35
   for c = 1:k
36
       g(:,c) = X'*(exp(XW(:,c))./Z-(y == c));
37
38
   end
   g = reshape(g, [p*k 1]);
```

```
40
   end
41
   function [w, f] = proxGradGroupL1(funObj, w, X, y, k, lambda, maxIter)
42
   % Minimize funOb(w) + lambda*sum(abs(w))
43
44
   % Evaluate initial objective and gradient of smooth part
45
46
   [f,g] = \text{funObj}(w,X,y,k,\text{lambda});
   funEvals = 1;
47
48
   L = 1;
49
   [n,p] = size(X);
50
   k = \max(y);
   while funEvals < maxIter
52
        % proximal-gradient step
54
        alpha = 1/L;
55
        W = reshape(w, [p k]);
56
        W_{new} = W;
57
        W_{update} = reshape(w - alpha*g, [p k]);
58
        for i = 1: p
59
             W_new(i,:) = softThreshold(W_update(i,:),alpha*lambda);
60
61
        end
        w_new = W_new(:);
62
        \%w_new = softThreshold(w - alpha*g, alpha*lambda);
63
        [f_{\text{new}}, g_{\text{new}}] = \text{funObj}(w_{\text{new}}, X, y, k, lambda);
64
        funEvals = funEvals + 1;
65
66
        % adaptive step-size
67
        while f_{\text{new}} > f + g'*(w_{\text{new}} - w) + (L/2)*norm(w_{\text{new}}-w)^2
             L = L*2;
69
             alpha = 1/L;
70
             w_new = softThreshold(w - alpha*g, alpha*lambda);
71
             [f_{\text{new}}, g_{\text{new}}] = \text{funObj}(w_{\text{new}}, X, y, k, lambda);
72
             funEvals = funEvals + 1;
73
        end
74
75
        w = w_n ew;
76
        f = f_new;
77
        g = g_new;
78
79
        % Print out how we are doing
80
        optCond = norm(w-softThreshold(w-g,lambda), 'inf');
81
        fprintf('%6d %15.5e %15.5e %15.5e\n', funEvals, alpha, f + lambda*sum(abs(w)
82
             ), optCond);
83
         if optCond < 1e-1
84
             break;
85
        end
86
   end
87
   end
88
89
   function [w] = softThreshold(w, threshold)
90
        \mathbf{w} = (\mathbf{w}./(\mathbf{norm}(\mathbf{w}))).*\mathbf{max}(0,\mathbf{norm}(\mathbf{w})-\mathbf{threshold});
91
92
   end
93
```

```
function [nll,g] = softmaxLoss(w,X,y,k,lambda)
95
    [n,p] = size(X);
96
   W = reshape(w, [p k]);
97
98
   XW = X*W:
99
   Z = sum(exp(XW), 2);
100
101
   ind = sub2ind([n k], [1:n]', y);
102
103
   tmp = sqrt(sum(W.^2,2));
104
105
   nll = -sum(XW(ind) - log(Z));
106
   g = zeros(p,k);
108
    for c = 1:k
109
        g(:,c) = X'*(exp(XW(:,c))./Z-(y == c));
110
111
   g = reshape(g, [p*k 1]);
112
   end
113
```

2.3 Stochastic Gradient

- 1. We tried the step-sizes defined by 1/(0.1t), 1/(0.2t)..., 1/(2t), and $1/(0.1\sqrt{t}), 1/(0.2\sqrt{t})..., 1/(2\sqrt{t})$. The maximum number of passes is set to 10. The step-size with the best performance is $\frac{1}{1.8\sqrt{t}}$ and its corresponding function value is 2.7145e+04.
- 2. We choose to weight the w with equal weights. The step-size with the best performance is $1/\sqrt{t}$ and its corresponding function value is 2.7076e+04.
- 3. We tried the step-sizes defined by $1/(0.1\sqrt{t}), 1/(0.2\sqrt{t})..., 1/(2\sqrt{t})$, and δ defined by 1,101,201,301,401. The sequence with the best performance is step-size = $1/(0.1\sqrt{t})$ and $\delta = 301$, the function value is 2.7083e+04. There are other combinations that gives close results: e.g. step-size = $1/(0.2\sqrt{t})$ and $\delta = 301$; step-size = $1/(0.3\sqrt{t})$ and $\delta = 301$.

```
1 load quantum.mat
  [n,d] = size(X);
  lambdaFull = 1;
3
4
  % Initialize
5
  maxPasses = 10;
   progTol = 1e-4;
  w = zeros(d,1);
  lambda = lambdaFull/n; % The regularization parameter on one example
  dt = zeros(d,d);
  tmp = zeros(d,1); % save the summation of the squared gradient in D
  % Stochastic gradient
  w_{-}old = w;
   c_{-candidate} = 0.1:0.1:2;
   for j = 1: length(c_candidate)
                                     %search for the constant value in the step
      sie
                            %search for the optimal delta
       for k = 1:100:500
16
```

```
w = zeros(d,1);
17
            w_{-}old = w;
18
            w_pre_iter = w;
19
            tmp = zeros(d,1);
20
            delta = k;
21
            c_candidate(j);
22
                 for t = 1: maxPasses*n
23
^{24}
                     % Choose variable to update
25
                     i = randi(n);
26
27
                     % Evaluate the gradient for example i
28
29
                     [f,g] = logisticL2\_loss(w,X(i,:),y(i),lambda);
                     tmp = tmp + g.^2;
31
                     D = diag(1./(sqrt(delta + tmp)));
32
33
                     % Choose the step-size
34
                     alpha = 1./(c_candidate(j)*t^(0.5));
35
36
                     w = w - alpha*D*g;
37
38
                     if mod(t,n) == 0
39
                         change = norm(w-w_old, inf);
40
                         %fprintf('Passes = %d, function = %.4e, change = %.4f\n',
41
                             t/n, logisticL2_loss(w,X,y,lambdaFull),change);
                          if change < progTol
42
                               fprintf('Parameters changed by less than progTol on
43
                             pass(n');
                              break;
44
                          end
45
                          w_{-}old = w;
46
                     end
                end
48
                  fprintf('c = %d, delta = %d, Passes = %d, function = %.4e \n',
49
                     c_c and idate (j), k, t/n, logistic L2_loss (w, X, y, lambda Full));
        end
50
51
  _{
m end}
```

4. The function value is 2.7068e+04 after 10 passes. It appears to converge faster than the algorithms in 1-3.

```
1 load quantum.mat
2 [n,d] = size(X);
3 lambdaFull = 1;
4
5 % Initialize
6 maxPasses = 10;
7 progTol = 1e-6;
8 w = zeros(d,1);
9 lambda = lambdaFull/n; % The regularization parameter on one example
10
11 % Stochastic gradient
12 w\_old = w;
```

```
13 L = .25*max(sum(X.^2, 2)) + lambda;
14
  D = zeros(d,1);
                      %represent the sum of the gradients calculated from n
15
      samples
   Y_i = zeros(n,d);
16
   g_{old} = zeros(d,1);
^{17}
18
   for t = 1: maxPasses*n
19
       % Choose variable to update
20
       i = randi(n);
^{21}
22
       % Evaluate the gradient for example i
23
24
       [f,g] = logisticL2\_loss(w,X(i,:),y(i),lambda);
25
26
       D = D - (Y_i(i, :))' + g;
27
       Y_{-i}(i, :) = g';
28
       \% Choose the step-size
30
       alpha = 1/L;
31
       w = w - (alpha*D./n);
32
33
       if \mod(t,n) == 0
34
            change = norm(w-w_old, inf);
35
            fprintf('Passes = %d, function = %.4e, change = %.4f\n',t/n,
36
               logisticL2_loss (w,X,y,lambdaFull),change);
            if change < progTol
37
                fprintf('Parameters changed by less than progTol on pass\n');
38
                break;
            end
40
            w_old = w;
41
       end
42
43
  end
```

3 Kernels and Duality

3.1 Fenchel Duality

1. Let $X = X, z = X\omega$.

$$\begin{split} f^*(u) &= \sup_{z} \left\{ u^T z - \frac{1}{2} \|z - y\|^2 \right\} \\ &= \sup_{z} \left\{ -\frac{1}{2} \|z - u - y\|^2 + \frac{1}{2} \|u\|^2 + u^T y \right\} \\ &= \frac{1}{2} \|u\|^2 + u^T y. \\ g^*(u) &= \sup_{\omega} \left\{ u^T \omega - \frac{\lambda}{2} \|\omega\|^2 \right\} \\ &= \sup_{\omega} \left\{ -\frac{\lambda}{2} \|\omega - \frac{1}{\lambda} u\|^2 + \frac{1}{2\lambda} \|u\|^2 \right\} \\ &= \frac{1}{2\lambda} \|u\|^2. \end{split}$$

Therefore, we have

$$D(z) = -f^*(-z) - g^*(X^T z) = -(\frac{1}{2} ||z||^2 - z^T y) - \frac{1}{2\lambda} ||X^T z||^2.$$

2. Let $X = X, z = X\omega$.

$$f^*(u) = \sup_{z} \left\{ u^T z - \|z - y\|_1 \right\}$$

$$= \sup_{z} \left\{ \sum_{i=1}^{n} \left(u_i z_i - |z_i - y_i| \right) \right\}$$

$$= \left\{ u^T y \quad \text{if } |u_i| \le 1 \text{ for all } i \right\}$$

$$= \left\{ u^T \omega - \lambda \|\omega\|_1 \right\}$$

$$= \left\{ 0 \quad \text{if } |u_i| \le \lambda \text{ for all } i \right\}$$

$$= \left\{ 0 \quad \text{if } |u_i| \le \lambda \text{ for all } i \right\}$$

$$= \left\{ 0 \quad \text{else} \right\}$$

Therefore, we have

$$D(z) = -f^*(-z) - g^*(X^T z) = \begin{cases} z^T y & \text{if } |z_i| \le 1 \text{ and } |(X^T z)_i| \le \lambda \text{ for all } i \\ -\infty & \text{else} \end{cases}.$$

3. Let
$$X = -\begin{pmatrix} y^1 x^1 \\ y^2 x^2 \\ \vdots \\ y^n x^n \end{pmatrix}$$
 and $z = X\omega$.
$$f^*(u) = \sup_{z} \left\{ u^T z - \sum_{i=1}^n \log(1 + \exp(z_i)) \right\}$$

$$= \begin{cases} \sum_{i=1}^n \left\{ u_i \log(u_i) + (1 - u_i) \log(1 - u_i) \right\} 1_{(0 < u_i < 1)} & \text{if } 0 \le u_i < 1 \text{ for all } i \\ \infty & \text{else} \end{cases}$$

$$g^*(u) = \sup_{\omega} \left\{ u^T \omega - \frac{\lambda}{2} \|\omega\|^2 \right\}$$

$$= \sup_{\omega} \left\{ -\frac{\lambda}{2} \|\omega - \frac{1}{\lambda} u\|^2 + \frac{1}{2\lambda} \|u\|^2 \right\}$$

$$= \frac{1}{2\lambda} \|u\|^2.$$

Therefore, we have

$$D(z) = -f^*(-z) - g^*(X^T z),$$

$$= \begin{cases} -\sum_{i=1}^n \left\{ -z_i \log(-z_i) + (1+z_i) \log(1+z_i) \right\} 1_{(-1 < z_i < 0)} - \frac{1}{2\lambda} ||X^T z||^2, & \text{if } -1 < z_i \le 0 \text{ for all } i \\ -\infty & \text{else} \end{cases}$$

3.2 Stochastic Dual Coordinate Ascent

The code is given below. The optimal values of the primal and dual objectives with $\lambda = 1$ are 90.6829 and 90.6827 respectively. The number of support vectors is 99.

```
%%%%%%%%%%%%%%%%%%%%%% Matlab code
X = load('statlog.heart.data');
y = X(:,end);
y(y==2) = -1;
X = X(:,1:end-1);
n = size(X,1);
% Add bias and standardize
X = [ones(n,1) standardizeCols(X)];
d = size(X,2);
% Set regularization parameter
lambda = 1;
% Some values used by the dual
```

```
YX = diag(y)*X;
G = YX*YX';

% Find min for -D(z)
z = quadprog(G/lambda, -ones(n,1), [],[],[], [], zeros(n,1), ones(n,1));
% Convert from dual to primal variables
w = (1/lambda)*(YX'*z);
% Evaluate primal objective:
P = sum(max(1-y.*(X*w),0)) + (lambda/2)*(w'*w)
% Evaluate dual objective:
D = sum(z) - (z'*G*z)/(2*lambda)
% count number of support vectors
sum(z> 0.01)
%%%%%%%%%%%%%%%% end-Matlab code
```

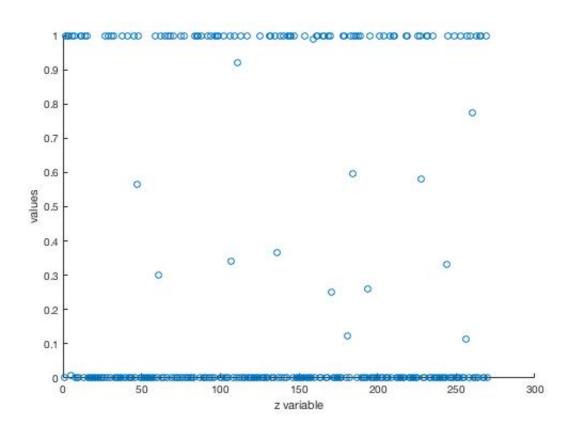


Figure 1: Values of z variables.

3.3 Large-Scale Kernel Methods

- 1. The best value for σ is 0.5 and the best value for λ is $\lambda_1 = 4.8828e 04$. The σ with Gaussian kernels is the same as the one with Gaussian RBFs. The λ for Gaussian RBFs was $\lambda_2 = 1.47e 05$. The approximate relationship between λ_1 and λ_2 is that $\lambda_2 \approx \lambda_1 * mean(K(:))$, where K = rbfBasis(X, X, sigma) is the matrix of pair-wise evaluations of the kernel for all training patterns.
- 2. The code is provided below and the performances with different m values (m=10,30,40,60,200) are visually examined by Figures 2-6. From the figures we see that the performance gets better when the m value increases. The performance of m=60 is as good as using the whole dataset (i.e. m=200). The squaredTestError values are 2.0787e+04 with m=10, 3.6422e+03 with m=30, 429.0070 with m=40, 82.8489 with m=60, and 83.4175 with m=200.

```
%%%%%%%%%%%%%%%%% Matlab code
lambda=4.8828e-04;
sigma = 0.5;
m=10; % or 30, 40, 60, 200
model = kernelRegression332(X, y, lambda, sigma, m);
yhat = model.predict(model, Xtest, m);
squaredTestError = sum((yhat-ytest).^2)/t
figure(1);
plot(X,y,'b.');
hold on
plot(Xtest, ytest, 'g.');
Xhat = [min(X):.1:max(X)]'; % Choose points to evaluate the function
yhat = model.predict(model, Xhat, m); % Gaussian kernel
plot(Xhat, yhat, 'r');
ylim([-300 400]);
function [model] = kernelRegression332(X,y,lambda,sigma, m)
  % Compute sizes
  [n,d] = size(X);
  % new
  K11 = rbfBasis(X(1:m,:),X(1:m,:),sigma);
  K21 = rbfBasis(X((m+1):n,:),X(1:m,:),sigma);
  K1 = [K11; K21];
  z = (K1' * K1 + lambda * K11) \ K1' * y;
  model.X = X;
  model.z = z;
  model.sigma = sigma;
```

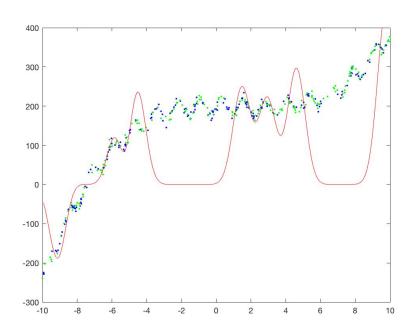


Figure 2: Performance with m = 10.

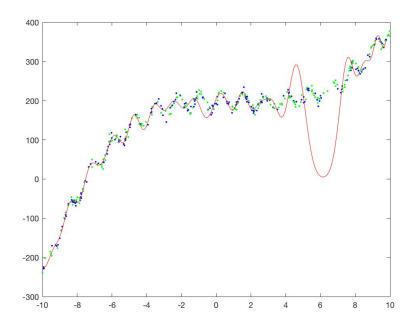


Figure 3: Performance with m=30.

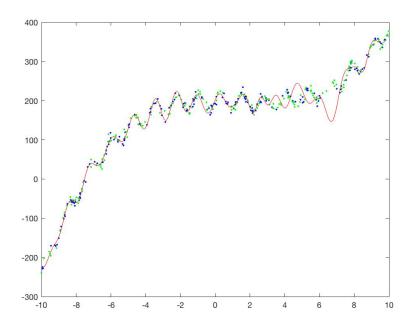


Figure 4: Performance with m=40.

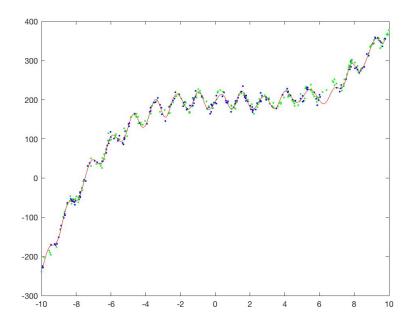


Figure 5: Performance with m = 60.

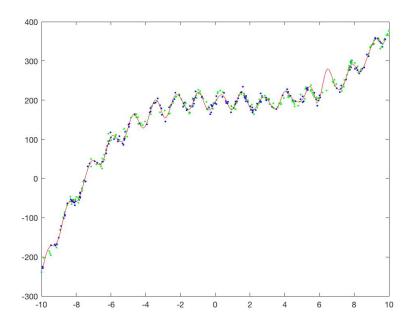


Figure 6: Performance with m = 200.

3. The code is given below and the performances with different m values (m = 3, 5, 8, 10, 20) are visually examined by Figures 7-11. From the figures we see that the performance gets better when the m value increases. This method can achieve a fitting curve that captures the overall pattern well with a small value of m (m=10), which is much smaller than the m value of the

subset of regressors model (m=60). However, this method does not capture the wiggly pattern in the data as the subset of regressors method does.

```
%%%%%%%%%%%%%%%% Matlab code
clear all
close all
% Load data
load nonLinear.mat % Loads {X,y,Xtest,ytest}
[n,d] = size(X);
[t, \tilde{}] = size(Xtest);
lambda=4.8828e-04;
sigma = 0.5;
rnq(1);
m=3; % or 5, 8, 10, 20
model = kernelRegression333(X, y, lambda, sigma, m);
yhat = model.predict(model, Xtest);
squaredTestError = sum((yhat-ytest).^2)/t
figure(1);
plot(X,y,'b.');
hold on
plot(Xtest, ytest, 'q.');
Xhat = [min(X):.1:max(X)]'; % Choose points to evaluate the function
yhat = model.predict(model, Xhat);
plot(Xhat, yhat, 'r');
ylim([-300 400]);
function [model] = kernelRegression333(X,y,lambda,sigma, m)
  % Compute sizes
  [n,d] = size(X);
  R = randn(d, m) *sigma; % dxm
  Z = \exp(\operatorname{sqrt}(-1) * X * R); % n, m
  % Solve least squares problem(
  z = (Z' * Z + lambda * eye(m)) \setminus Z' * y; % m, 1
  model.R = R;
                 % dxm
  model.z = z;
                 % mx1
  model.sigma = sigma;
  model.predict = @predict;
end
```

```
function [yhat] = predict(model, Xhat)
  Zhat = exp(sqrt(-1)*Xhat*model.R); % n2xm
  yhat = Zhat*model.z;
end
%%%%%%%%%%%%%%%%%%% end- Matlab code
```

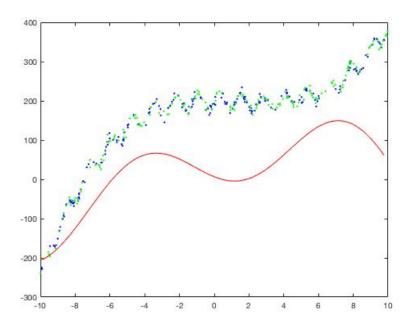


Figure 7: Performance with m = 3.

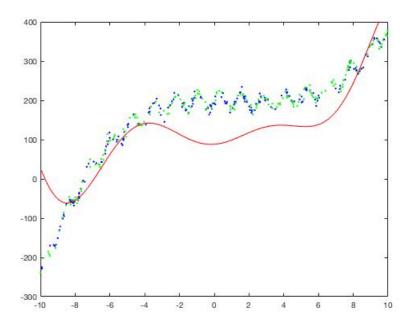


Figure 8: Performance with m=5.

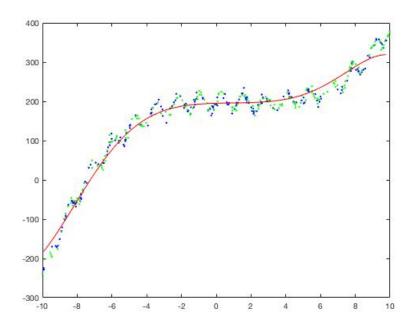


Figure 9: Performance with m = 8.

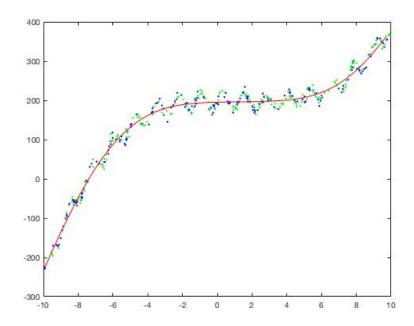


Figure 10: Performance with m = 10.

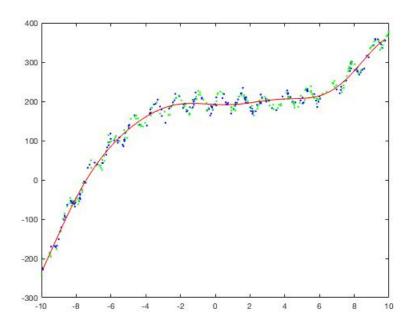


Figure 11: Performance with m = 20.