LAGraph Algorithms

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Abstract

Theoretical documentation for LAGraph.

1 Notation

Table 1 contains the notation of GraphBLAS operations

Additionally, we use D = diag(J, n) to construct a diagonal matrix $D \leftarrow \{J, J, [1, 1, ..., 1]\}$. The elements of the matrix are D(j, j) = 1 for $j \in J$.

Initializing scalars, vectors, and matrices (GraphBLAS methods):

- s = fp64()
- $\mathbf{u} = fp32(\mathbf{n})$
- A = uint16(m, n)
- A = int64(k, m)

2 Algorithms

LAGraph [12] implements graph algorithms using the GraphBLAS C API [13]. An incomplete list of GraphBLAS algorithms:

- MSBFS [8], bidirectional BFS [8], pushpull BFS [19]
- DFS [16]
- weakly connected components [21]
- SCC (LAGraph)
- SSSP, delta-stepping [17]
- triangle count [1, 5], triangle enumeration [1], item local clustering coefficient [4]
- *k*-truss [5]
- betweenness centrality [12]
- closeness centrality [8]
- DNN algorithm [10, 7]
- PageRank variants (at least 2) [4], IISWC paper, ...
- Louvain [11]
- property graphs: incremental TTC case [9], SIGMOD 2014 Contest [8] Roi Lipman's talk¹
- CFPQs based on a string of GRADES/ADBIS/other papers [2, 3, 14, 18, 15]
- Implementations: GBTL², SuiteSparse:GraphBLAS [6], GraphBLAST [20]

 $^{^{1}} http://wiki.ldb council.org/pages/viewpage.action?pageId=106233859 \& preview=/106233859/111706128/LDBC-July-2019.pdf$

²https://github.com/cmu-sei/gbtl

| op./method | name | notation |
|-------------------|---|--|
| mxm vxm mxv | matrix-matrix multiplication vector-matrix multiplication matrix-vector multiplication | $C < M > + = A + . \times B$ $w < m > + = u + . \times A$ $w < m > + = A + . \times u$ |
| eWiseAdd | element-wise addition set union of patterns | C < M > + = A + B $w < m > + = u + v$ |
| eWiseMult | element-wise multiplication set intersection of patterns | $C < M > + = A \times B$ $w < m > + = u \times v$ |
| extract | extract submatrix extract column vector extract row vector extract subvector | |
| assign | assign matrix to submatrix with mask for C assign scalar to submatrix with mask for C assign vector to subvector with mask for W assign scalar to subvector with mask for W | |
| subassign (GxB) | assign matrix to submatrix with submask for $\mathtt{C}(\mathtt{I},\mathtt{J})$ assign scalar to submatrix with submask for $\mathtt{C}(\mathtt{I},\mathtt{J})$ assign vector to subvector with submask for $\mathtt{w}(\mathtt{I})$ assign scalar to subvector with submask for $\mathtt{w}(\mathtt{I})$ | C(I, J) < M>+ = A C(I, J) < M>+ = s w(I) < m>+ = u w(I) < m>+ = s |
| apply | apply unary operator | $ \begin{array}{l} \mathtt{C} < \mathtt{M} > + = \mathbf{f}(\mathtt{A}) \\ \mathtt{w} < \mathtt{m} > + = \mathbf{f}(\mathtt{u}) \end{array} $ |
| select (GxB) | apply select operator | $ \begin{split} & \texttt{C} < \texttt{M} > + = select(\texttt{A}, \textbf{f}(\texttt{k})) \\ & \texttt{C} < \texttt{M} > + = select(\texttt{low} \leq \texttt{A} \leq \texttt{up}) \\ & \texttt{w} < \texttt{m} > + = select(\texttt{u}, \textbf{f}(\texttt{k})) \\ & \texttt{w} < \texttt{m} > + = select(\texttt{low} \leq \texttt{u} \leq \texttt{up}) \end{split} $ |
| reduce | reduce matrix to column vector reduce matrix to scalar reduce vector to scalar | w < m > + = [+ A] s + = [+ A] s + = [+ u] |
| transpose | transpose | C < M > + = A' |
| kronecker | Kronecker multiplication | C < M > + = kron(A, B) |
| new | new matrix new vector | $\begin{aligned} & \overset{\textbf{A}}{\textbf{u}} = TYPEPRECISION(\textbf{n}, \textbf{m}) \\ & \overset{\textbf{u}}{\textbf{u}} = TYPEPRECISION(\textbf{n}) \end{aligned}$ |
| build | build matrix from index/value arrays build vector from index/value arrays | $ \begin{array}{l} \textbf{C} < - \{ \texttt{I}, \texttt{J}, \texttt{X} \} \\ \textbf{w} < - \{ \texttt{I}, \texttt{X} \} \end{array} $ |
| extractTuples | extract index/value arrays | {I, J, X}<- A {I, X}<- u |
| dup | duplicate matrix duplicate vector | C<- A w<- u |
| extractElement | extract scalar element | s = A(i,j) $s = u(i)$ |
| setElement | set element | $egin{aligned} \mathbf{C}(\mathtt{i},\mathtt{j}) &= \mathtt{s} \\ \mathtt{w}(\mathtt{i}) &= \mathtt{s} \end{aligned}$ |

Table 1: GraphBLAS operations and methods based on [6]. Notation: Matrices and vectors are typeset in bold, starting with uppercase (A) and lowercase (u) letters, respectively. Scalars including indices are lowercase italic (s, i, j) while arrays are uppercase italic (X, I, J). + and × are the addition and multiplication operators forming a semiring and default to conventional arithmetic + and × operators. \odot is the apply operator. Masks <M> and <m> are used to selectively write to the result matrix/vector. The complements of masks <M>, <m> can be selected with the negation symbol, denoted with <!M>, <!m>, respectively. Masks with "replace" semantics (annihilating all elements outside the mask) are denoted with <<M>>>, <<!M>>>, <<m>>>, and <<!m>>>, The structure of the mask is denoted with <{M}>, <!m}>, and <!{m}>>, and <!{m}>>.

Algorithm 1: Breadth-first search.

Algorithm 2: Breadth-first search (push/pull).

```
Function BFS
frontier(s) = TRUE
for level = 1 to n - 1 do

seen < frontier >= level
push = use some heuristics to determine whether to push/pull
frontier <<!seen >>= frontier any.pair A

selse
frontier <<!seen >>= A any.pair frontier
```

Algorithm 3: Multi-source breadth-first search.

```
Data: ...
Result: ...

Function MSBFS

Frontier <- diag(S,n)

for level = 1 to n - 1 do

Seen <Frontier >= level

Frontier <<!Seen >>= Frontier any.pair A
```

Algorithm 4: Betweenness centrality.

```
1 Function MSBFS
      // The NumSp structure holds the number of shortest paths for each node and starting
         vertex discovered so far.
      // Initialized to source vertices.
3
      NumSp < -\{s, [1, 1, ..., 1]\}
4
      // The Frontier holds the number of shortest paths for each node and starting vertex
5
         discovered so far.
      // Initialized to source vertices.
6
      Frontier < NumSp>= A(s,:)
7
      d = 0
8
      // The Sigmas matrices store frontier information for each level of the BFS phase.
9
      // BFS phase (forward sweep)
10
11
      do
         // Sigmas[d](:,s) = d^{th} level frontier from source vertex s
12
         Sigmas[d] = bool(n, nsver)
13
         Sigmas[d](:,:) = Frontier
                                                                     // Convert matrix to Boolean
14
         NumSp = NumSp + Frontier
                                                                        // Accumulate path counts
15
        Frontier << NumSp>>= A' + . × Frontier
                                                                                // Update frontier
16
      while nvals(Frontier) > 0
17
      NumSpInv = fp32(n, nsver)
18
      NumSpInv = 1.0DIV NumSp
19
      BCU = fp32(n, nsver)
20
      BCU(:) = 1.0
                                              // Make BCU dense, initialize all elements to 1.0
21
      W = fp32(n, nsver)
22
      // Tally phase (backward sweep)
23
      for i = d - 1 downto 0 do
24
         W << Sigmas[i] >>= NumSpInv DIV BCU
25
         \mathbb{W} \ll \operatorname{Sigmas}[i-1] >>= \mathbb{A} + . \times \mathbb{W} // Add contributions by successors and mask with that
26
          BFS level's frontier.
        BCU += W \times NumSp
27
      // Row reduce BCU and subtract nsver from every entry to account for 1 extra value
28
         per BCU row element
      delta = [+BCU]
29
      delta MINUS = nsver
30
```

Algorithm 5: PageRank (used in Graphalytics).

```
Data: alpha constant (damping factor)
  Result: ...
1 Function PageRank
     pr(:) = 1/n
      outdegrees = [+_i A(:,j)]
3
     for k = 1 to numIterations do
4
         importance = pr DIV outdegrees
5
         importance = times(importance, alpha)
                                                           // apply the times(x,s) = x \cdot s operator
6
         importance = importance + . \times A
         danglingVertexRanks < !outdegrees >= pr(:)
8
         totalDanglingRank = [+ danglingVertexRanks(i)](alpha)/(n)
9
         pr(:) = (1 - alpha)/(n) + totalDanglingRank
10
         pr < importance >= pr + importance
11
```

Algorithm 6: Algebraic Bellman-Ford for SSSP.

Algorithm 7: Delta-stepping SSSP.

```
Data:
          A, A_H, A_L \in fp(|V|, |V|)
          \mathtt{s},\mathtt{i} \in \mathtt{uint}()
          \Delta \in \mathtt{fp}()
          \mathtt{t},\mathtt{t}_{\mathtt{Req}}\in\mathtt{fp}(|\mathtt{V}|)
          \mathtt{t}_{B_{\mathtt{i}}}, \mathtt{e} \in \mathtt{uint}(|\mathtt{V}|)
  1 Function DeltaStepping
            A_L = select(0 < A \le \Delta)
            {\bf A_H} = select(\Delta < {\bf A})
  3
  4
            t(:) = \infty
            t(s) = 0
  5
            while nvals(select(i\Delta \leq t)) \neq 0 do
  6
                   \mathtt{t}_{\mathtt{B}_{\mathtt{i}}} = select(\mathtt{i}\Delta \leq \mathtt{t} < (\mathtt{i}+1)\Delta)
  8
                   while t_{B_i} \neq 0 do
  9
                          t_{Req} = A_L \, ' + . \times (t \times t_{B_i})
10
                          e = select(0 < e + t_{B_i})
11
                          \mathtt{t_{B_i}} = \mathit{select}(\mathtt{i}\Delta \leq \mathtt{t_{Req}} < (\mathtt{i}+1)\Delta) \times (\mathtt{t_{Req}} \min_{+} \mathtt{t})
12
                   t_{Req} = A_H' + . \times (t \times e)
13
                   t = t \min t_{Req}
14
                   i = i + 1
15
```

Algorithm 8: All-pairs shortest path (Floyd-Warshall algorithm).

Algorithm 9: FastSV algorithm.

```
1 Function FastSV
 \mathbf{2}
       n = nrows(A)
       gf = f
 3
       dup = gf
 4
       mngf = gf
 5
       \{I,X\} \leftarrow f
 6
       repeat
 7
          // Step 1: Stochastic hooking
 8
           mngf = mngf min A
 9
           mngf = mngf second.min gf
10
           f(X) = f \min mngf
11
           // Step 2: Aggressive hooking
12
13
           {\tt f} = {\tt fminmngf}
14
           // Step 3: Shortcutting
           f = f \min gf
15
           // Step 4: Calculate grandparents
16
           \{I,X\} \leftarrow f
17
           gf = f(X)
18
           // Step 5: Check termination
19
           diff = dup \neq gf
20
           sum = [+_i diff(i)]
21
           dup = gf
\mathbf{22}
       \mathbf{until}\ \mathbf{sum} == 0
23
```

Algorithm 10: Triangle count (Cohen's algorithm).

Algorithm 11: Triangle count (Sandia).

Algorithm 12: Triangle count (FLAME).

Algorithm 13: Local clustering coefficient.

Algorithm 14: k-truss algorithm.

```
1 Function KTruss
     C <- A
2
     nonzeros < - nvals(C)
3
     for i = 1 to n - 1 do
4
         C < C > = C + . land C
5
         C = select(C \ge k - 2)
6
         if nonzeros == nvals(C) then
7
          break
8
9
        nonzeros \leftarrow nvals(C)
```

Algorithm 15: Louvain algorithm (WIP).

```
1 Function Louvain
      G += G'
 \mathbf{2}
      k = [+A]
 3
      m = (1)/(2)[+k]
 4
      S <- I
 5
 6
      vertices_changed <- nvals(k)</pre>
 7
      while vertices_changed > 0 do
 8
          for j \in range(|V|) do
 9
10
              v = G(j,:)
              t_q = v any.pair S
11
              sr = S(j,:)
12
              S(j,:) = empty
13
              q <- k
15
              q < k > x = -k(j)/m
16
              q + = v
17
18
              q_1 < t_q >= q + . \times S
19
              t = (q_1 == [\max q_1])
20
              while nvals(t) \neq 1 do
21
                 p = random() \times t
22
                t = (p == [\max p])
23
              S(j,:) = t
24
25
              if nvals(sr \times t) == 0 then
26
               vertices\_changed = nvals(k)
27
              vertices\_changed = vertices\_changed - 1
28
```

Algorithm 16: Community detection using label propagation (for undirected graphs).

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