

$$A = [a_{ij}] = \begin{bmatrix} a_{11} & a_{12} & a_{13} & \cdots & a_{1n} \\ a_{21} & a_{22} & a_{23} & \cdots & a_{2n} \\ a_{31} & a_{32} & a_{33} & \cdots & a_{3n} \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ a_{m1} & a_{m2} & a_{m3} & \cdots & a_{mn} \end{bmatrix}$$

$$\begin{aligned} A^2 XBA &= AB \\ A^{-2} A^2 XBA &= A^{-2} AB \\ XBA &= A^{-1} B \\ XBAA^{-1} &= A^{-1} BA^{-1} \\ XB &= A^{-1} BA^{-1} \\ XBB^{-1} &= A^{-1} BA^{-1} B^{-1} \\ X &= A^{-1} BA^{-1} B^{-1} \end{aligned}$$

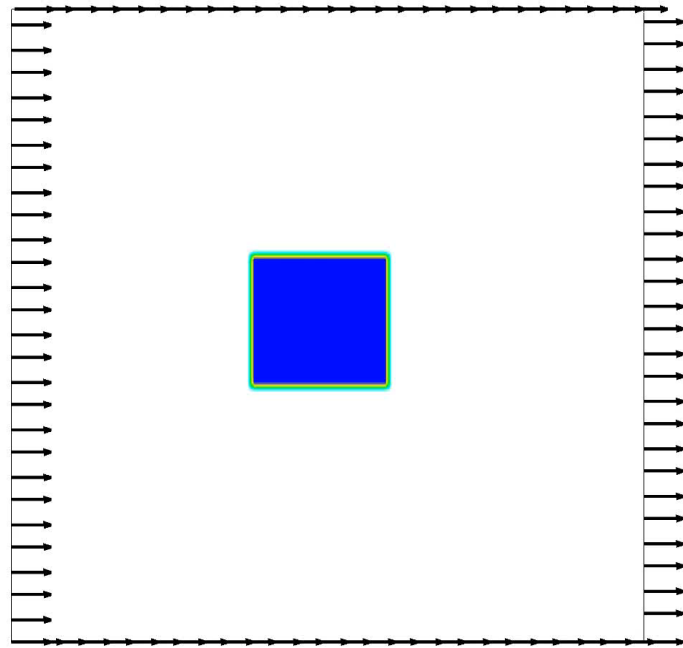
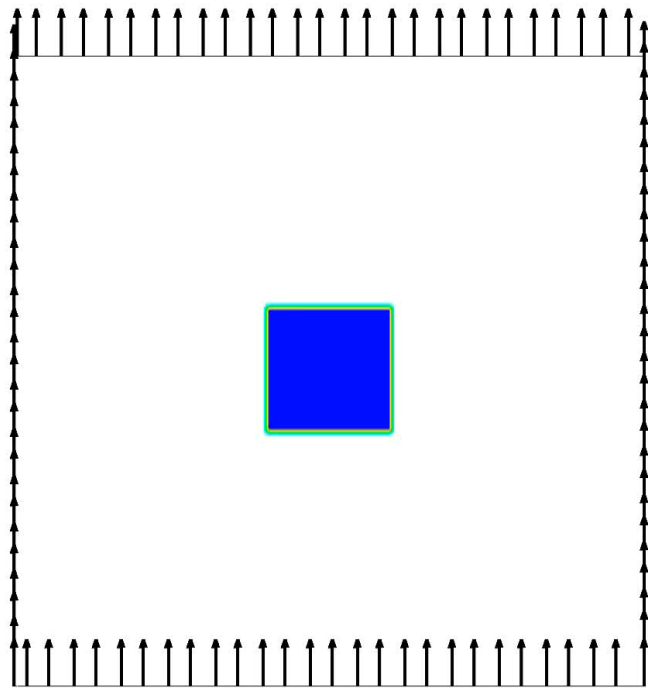
Stabilization of Gauss-Seidel Smoother

NAOFE, Dong-A university

Yeoun Joo KIM

Motivation

Laplace-Trans
$$F(p) = \int_0^{\infty} e^{-pt} f(t) dt,$$

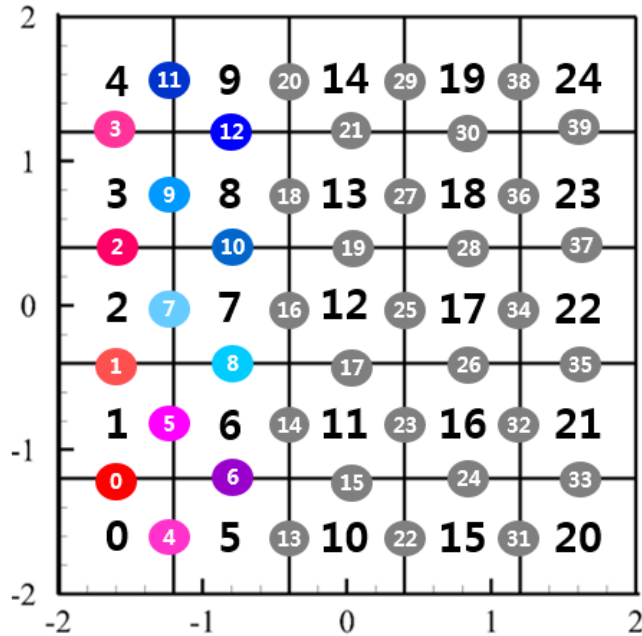


➤ x방향으로 흐르지 못하는 이유는?

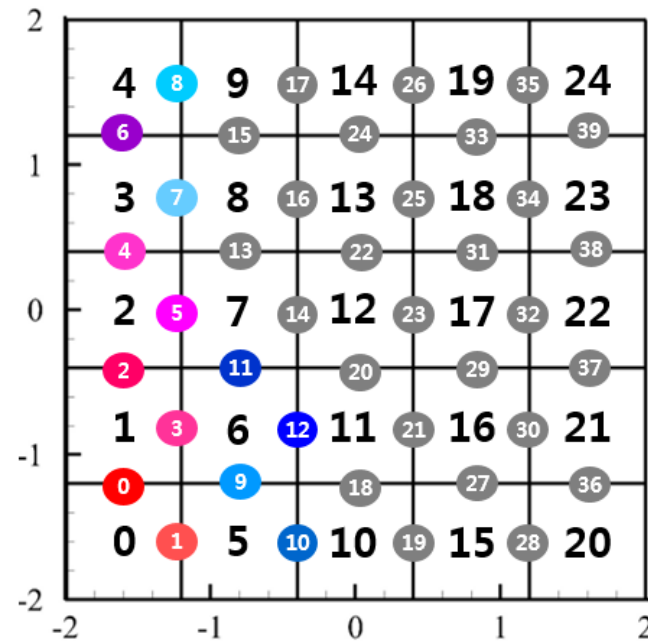
Cases of Face-Ordering

Laplace-Transform

$$F(p) = \int_0^{\infty} e^{-pt} f(t) dt$$



Grid 1



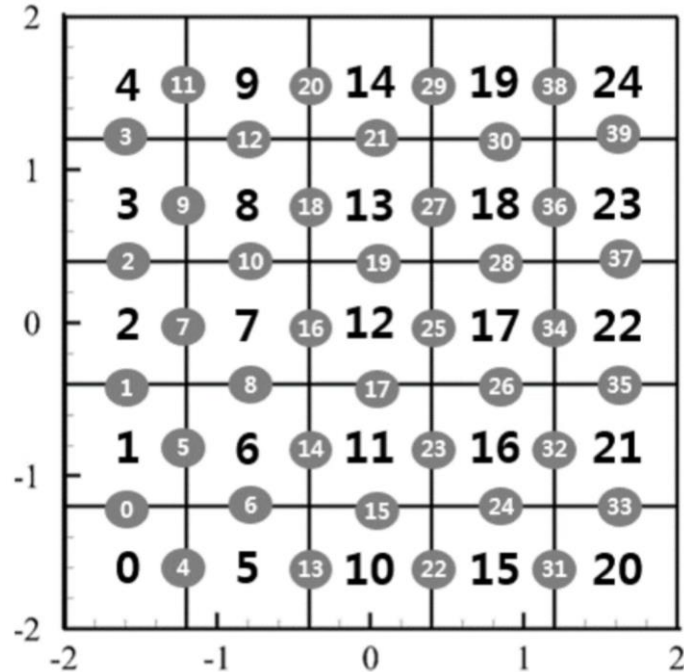
Grid 2

Different ordering of face index in two grid systems

GS Smoother : Grid 1

Laplace-Trans

$$F(p) = \int_0^{\infty} e^{-pt} f(t) dt$$



```
for (label celli=0; celli<nCells; celli++)
{
    // Start and end of this row
    fStart = fEnd;
    fEnd = ownStartPtr[celli + 1];

    // Get the accumulated neighbour side
    psii = bPrimePtr[celli];

    // Accumulate the owner product side
    for (label facei=fStart; facei<fEnd; facei++)
    {
        psii -= upperPtr[facei]*psiPtr[uPtr[facei]];
    }
}
```

$$\varphi_{i,j} = l_{i-1}\varphi_{i-1,j} + u_{i+1}\varphi_{i+1,j} \quad \text{missed}$$

$$+ l_{j-1}\varphi_{i,j-1} + u_{j+1}\varphi_{i,j+1}$$

$$+ d_{k,l}\varphi_{k,l}$$

$$\begin{cases} k \neq i-1, i+1 \\ l \neq j-1, j+1 \end{cases}$$

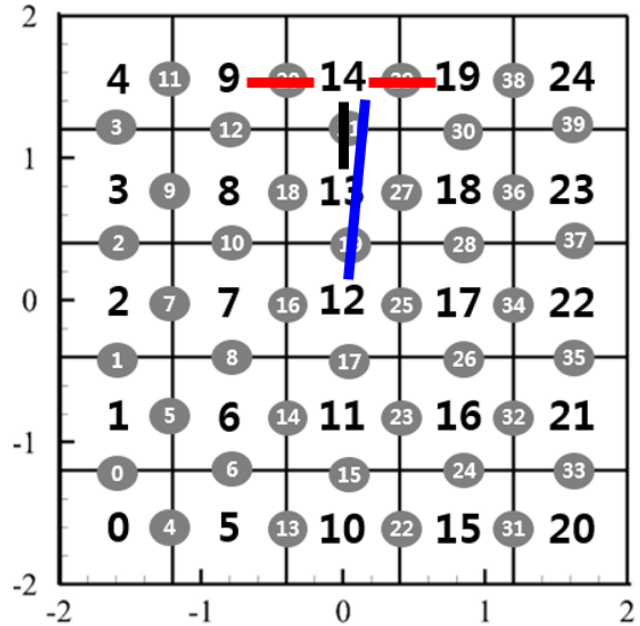
misplaced

GS Smoother : Grid 1

Laplace-Transform

$$F(p) = \int_0^{\infty} e^{-pt} f(t) dt$$

	0	1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	16	17	18	19	20	21	22	23	24		
0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
1	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	1
2	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	2
3	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	3
4	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	4
5	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	5
6	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	6
7	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	7
8	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	8
9	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	9
10	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	10
11	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	11
12	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	12
13	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	13
14	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	14
15	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	15
16	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	16
17	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	17
18	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	18
19	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	19
20	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.2	0	0	0	0	0	20
21	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.2	0	0	0	21
22	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.2	0	0	22
23	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.2	0	23
24	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.2	0	24

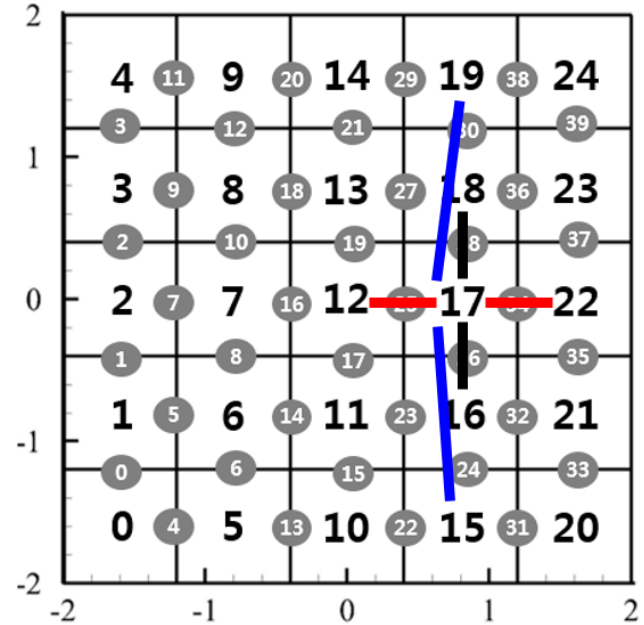


GS Smoother : Grid 1

Laplace-Trans

$$F(p) = \int_0^{\infty} e^{-pt} f(t) dt$$

	0	1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	16	17	18	19	20	21	22	23	24		
0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
1	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	1
2	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	2
3	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	3
4	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	4
5	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	5
6	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	6
7	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	7
8	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	8
9	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	9
10	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	10
11	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	11
12	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0.997506	3.2	12
13	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	13
14	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0.0024876	0	14
15	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	15
16	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	16
17	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	17
18	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	18
19	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	19
20	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.2	0	0	0	0	0	20
21	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.2	0	0	0	21
22	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	0	3.2	0	22
23	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	0	0	23
24	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.2	0	24

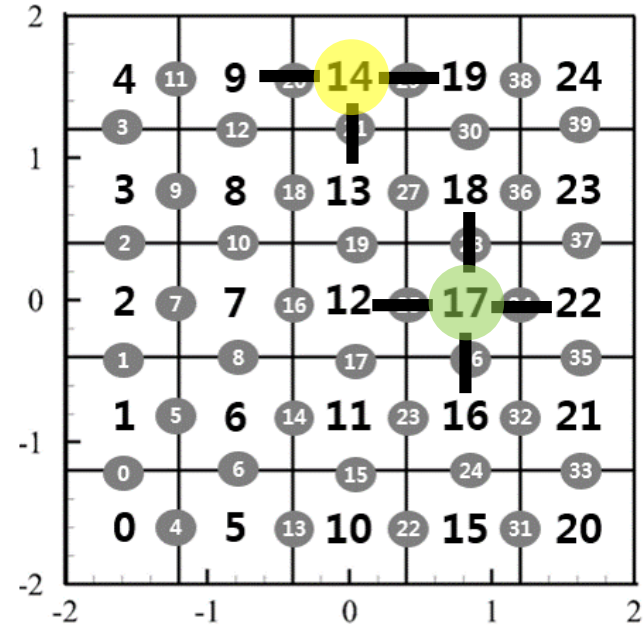


GS Smoother : Grid 1

Laplace-Trans

$$F(p) = \int_0^{\infty} e^{-pt} f(t) dt$$

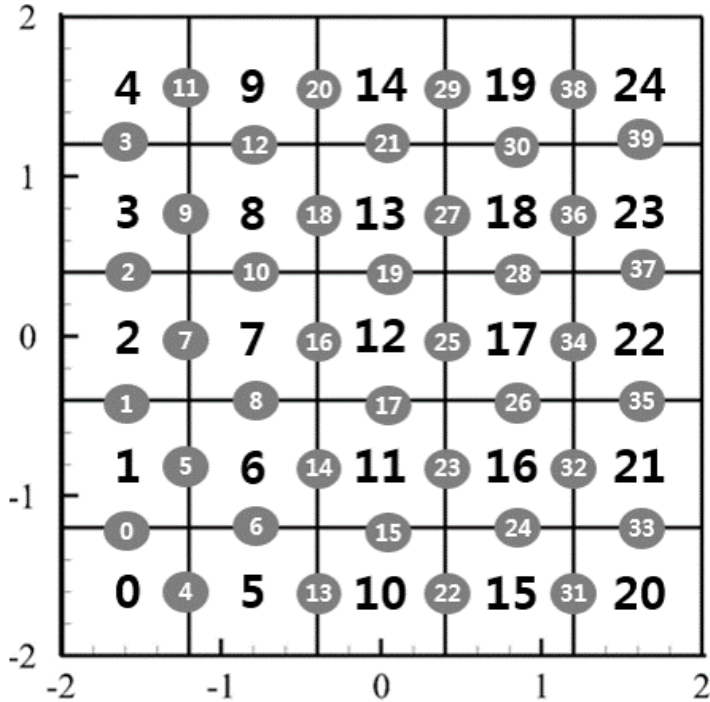
	0	1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	16	17	18	19	20	21	22	23	24		
0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
1	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	1
2	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	2
3	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	3
4	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	4
5	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	5
6	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	6
7	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	7
8	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	8
9	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	9
10	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	10
11	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	11
12	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	12
13	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	13
14	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	14
15	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	15
16	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.208	0	0	0	0	0	0	0	0	16
17	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.208	0	0	0	0	0	0	0	17
18	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.208	0	0	0	0	0	0	18
19	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.208	0	0	0	0	0	19
20	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.2	0	0	0	0	20
21	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.2	0	0	0	21
22	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.2	0	0	22
23	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.2	0	23
24	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.2	0	24



DILU Smoother : Grid 1

Laplace-Transform

$$F(p) = \int_0^{\infty} e^{-pt} f(t) dt$$



```
for (label face=0; face<nFaces; face++)
{
    rAPtr[uPtr[face]] -= lowerPtr[face]*psiPtr[lPtr[face]];
    rAPtr[lPtr[face]] -= upperPtr[face]*psiPtr[uPtr[face]];
}

rA *= rD_;

label nFaces = matrix_.upper().size();
for (label face=0; face<nFaces; face++)
{
    label u = uPtr[face];
    rAPtr[u] -= rDPtr[u]*lowerPtr[face]*rAPtr[lPtr[face]];
}
```

```
label nFacesM1 = nFaces - 1;
for (label face=nFacesM1; face>=0; face--)
{
    label l = lPtr[face];
    rAPtr[l] -= rDPtr[l]*upperPtr[face]*rAPtr[uPtr[face]];
}

psi += rA;
```

$$Ax_i = B_i$$

$$A(x_i - x_{i-1}) = B_i - B_{i-1}$$

$$(D + L + U)(x_i - x_{i-1}) = B_i - B_{i-1}$$

$$x_i - x_{i-1} \cong D^{-1}(B_i - B_{i-1})$$

$$x_i - x_{i-1} + D^{-1}L(x_i - x_{i-1}) + D^{-1}U(x_i - x_{i-1}) = D^{-1}(B_i - B_{i-1})$$

$$x_i \cong x_{i-1} + D^{-1}(B_i - B_{i-1}) - D^{-1}LD^{-1}(B_i - B_{i-1}) - D^{-1}UD^{-1}(B_i - B_{i-1})$$

$$rA = D^{-1}(B_i - B_{i-1})$$

$$x_i \cong x_{i-1} + rA - D^{-1}LrA - D^{-1}UrA$$

$$\|D(x_i - x_{i-1})\| \gg \|L(x_i - x_{i-1})\|$$

$$\|D(x_i - x_{i-1})\| \gg \|U(x_i - x_{i-1})\|$$

$rAPtr[u] -= rDPtr[u]*lowerPtr[face]*rAPtr[lPtr[face]];$

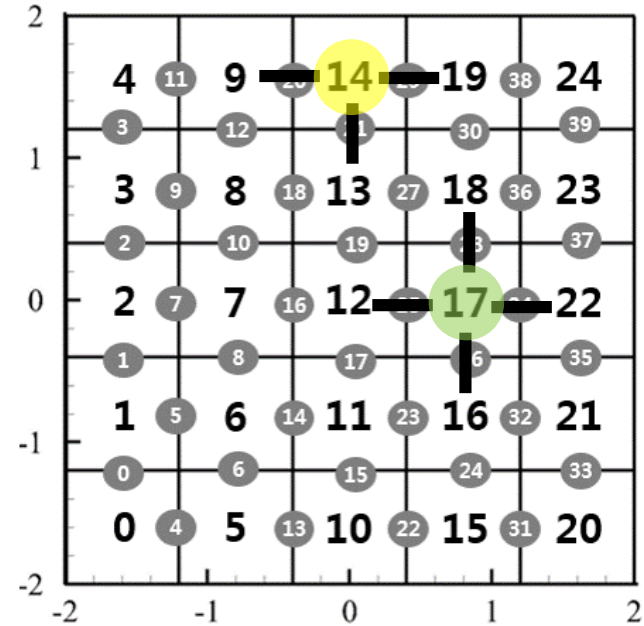
$rAPtr[l] -= rDPtr[l]*upperPtr[face]*rAPtr[uPtr[face]];$

DILU Smoother : Grid 1

Laplace-Trans

$$F(p) = \int_0^{\infty} e^{-pt} f(t) dt$$

	0	1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	16	17	18	19	20	21	22	23	24		
0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
1	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	1
2	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	2
3	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	3
4	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	4
5	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	5
6	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	6
7	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	7
8	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	8
9	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	9
10	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	10
11	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	11
12	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	12
13	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	13
14	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	14
15	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	15
16	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.208	0	0	0	0	0	0	0	0	16
17	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.208	0	0	0	0	0	0	0	17
18	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.208	0	0	0	0	0	0	18
19	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.208	0	0	0	0	0	19
20	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.2	0	0	0	0	20
21	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	0	3.2	0	0	21
22	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	0	3.2	0	22
23	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	0	3.2	23
24	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.2	0	24



Other Smoothers

Laplace-Trans
$$F(p) = \int_0^{\infty} e^{-pt} f(t) dt,$$

DILUGaussSeidel solver

```
diluSmoother_.smooth(psi, source, cmpt, nSweeps);  
gsSmoother_.smooth(psi, source, cmpt, nSweeps);
```

symGaussSeidel solver

```
for (label celli=0; celli<nCells; celli++)  
{  
    // Start and end of this row  
    fStart = fEnd;  
    fEnd = ownStartPtr[celli + 1];  
  
    // Get the accumulated neighbour side  
    psii = bPrimePtr[celli];  
  
    // Accumulate the owner product side  
    for (label facei=fStart; facei<fEnd; facei++)  
    {  
        psii -= upperPtr[facei]*psiPtr[uPtr[facei]];  
    }  
  
    // Finish current psi  
    psii /= diagPtr[celli];  
  
    // Distribute the neighbour side using current psi  
    for (label facei=fStart; facei<fEnd; facei++)  
    {  
        bPrimePtr[uPtr[facei]] -= lowerPtr[facei]*psii;  
    }  
  
    psiPtr[celli] = psii;  
}  
  
fStart = ownStartPtr[nCells];  
  
fStart = ownStartPtr[nCells];  
{  
    // Start and end of this row  
    fEnd = fStart;  
    fStart = ownStartPtr[celli];  
  
    // Get the accumulated neighbour side  
    psii = bPrimePtr[celli];  
  
    // Accumulate the owner product side  
    for (label facei=fStart; facei<fEnd; facei++)  
    {  
        psii -= upperPtr[facei]*psiPtr[uPtr[facei]];  
    }  
  
    // Finish psi for this cell  
    psii /= diagPtr[celli];  
  
    // Distribute the neighbour side using psi for this cell  
    for (label facei=fStart; facei<fEnd; facei++)  
    {  
        bPrimePtr[uPtr[facei]] -= lowerPtr[facei]*psii;  
    }  
  
    psiPtr[celli] = psii;  
}
```

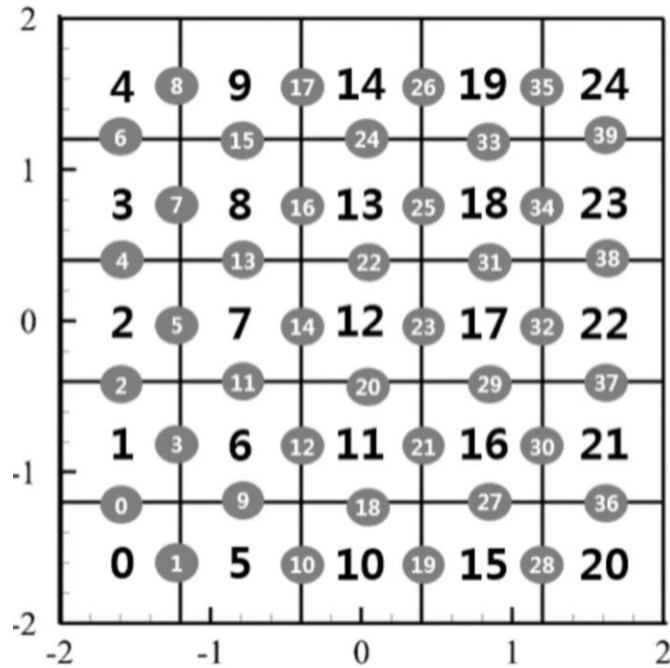
nonBlockingGaussSeidel solver

```
for (label celli=0; celli<blockStart; celli++)  
{  
    // Start and end of this row  
    fStart = fEnd;  
    fEnd = ownStartPtr[celli + 1];  
  
    // Get the accumulated neighbour side  
    curPsi = bPrimePtr[celli];  
  
    // Accumulate the owner product side  
    for (label curFace=fStart; curFace<fEnd; curFace++)  
    {  
        curPsi -= upperPtr[curFace]*psiPtr[uPtr[curFace]];  
    }  
  
    // Finish current psi  
    curPsi /= diagPtr[celli];  
  
    // Distribute the neighbour side using current psi  
    for (label curFace=fStart; curFace<fEnd; curFace++)  
    {  
        bPrimePtr[uPtr[curFace]] -= lowerPtr[curFace]*curPsi;  
    }  
  
    psiPtr[celli] = curPsi;  
}  
  
for (label celli=blockStart; celli < nCells; celli++)  
{  
    // Start and end of this row  
    fStart = fEnd;  
    fEnd = ownStartPtr[celli + 1];  
  
    // Get the accumulated neighbour side  
    curPsi = bPrimePtr[celli];  
  
    // Accumulate the owner product side  
    for (label curFace=fStart; curFace<fEnd; curFace++)  
    {  
        curPsi -= upperPtr[curFace]*psiPtr[uPtr[curFace]];  
    }  
  
    // Finish current psi  
    curPsi /= diagPtr[celli];  
  
    // Distribute the neighbour side using current psi  
    for (label curFace=fStart; curFace<fEnd; curFace++)  
    {  
        bPrimePtr[uPtr[curFace]] -= lowerPtr[curFace]*curPsi;  
    }  
  
    psiPtr[celli] = curPsi;  
}
```

➤ Based on GaussSeidel solver

GS Smoother : Grid 2

Laplace-Transform
$$F(p) = \int_0^{\infty} e^{-pt} f(t) dt$$



```
for (label celli=0; celli<nCells; celli++)  
{  
    // Start and end of this row  
    fStart = fEnd;  
    fEnd = ownStartPtr[celli + 1];  
  
    // Get the accumulated neighbour side  
    psii = bPrimePtr[celli];  
  
    // Accumulate the owner product side  
    for (label facei=fStart; facei<fEnd; facei++)  
    {  
        psii -= upperPtr[facei]*psiPtr[uPtr[facei]];  
    }  
}
```

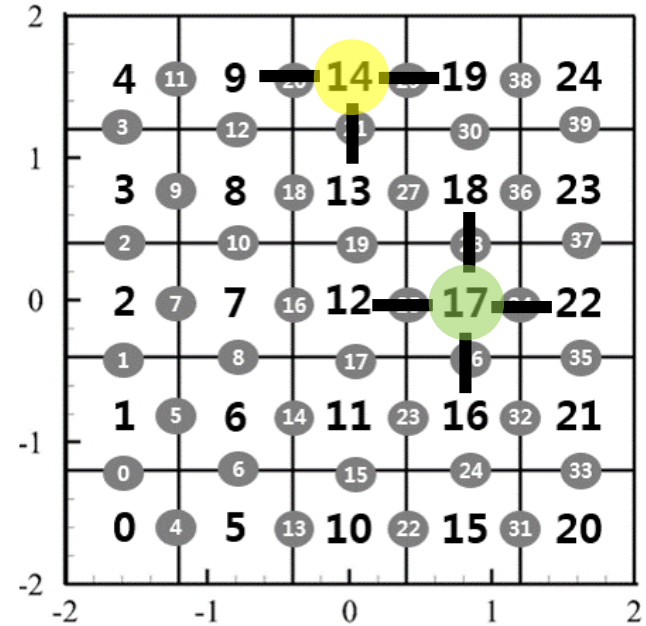
$$\begin{aligned}\varphi_{i,j} = & l_{i-1}\varphi_{i-1,j} + u_{i+1}\varphi_{i+1,j} \\ & + l_{j-1}\varphi_{i,j-1} + u_{j+1}\varphi_{i,j+1}\end{aligned}$$

GS Smoother : Grid 2

Laplace-Trans

$$F(p) = \int_0^{\infty} e^{-pt} f(t) dt$$

	0	1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	16	17	18	19	20	21	22	23	24		
0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
1	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	1
2	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	2
3	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	3
4	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	4
5	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	5
6	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	6
7	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	7
8	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	8
9	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	9
10	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	10
11	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	11
12	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	12
13	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	13
14	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	14
15	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	15
16	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.208	0	0	0	0	0	0	0	0	16
17	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.208	0	0	0	0	0	0	0	17
18	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.208	0	0	0	0	0	0	18
19	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.208	0	0	0	0	0	19
20	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.2	0	0	0	0	20
21	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	0	3.2	0	0	21
22	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	0	3.2	0	22
23	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	0	3.2	23
24	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.2	0	24



Code Analysis of GS

Laplace-Transform

$$F(p) = \int_0^{\infty} e^{-pt} f(t) dt$$

```

scalar psii;
label fStart;
label fEnd = ownStartPtr[0]; → start owner face at each cell

for (label celli=0; celli<nCells; celli++)
{
    // Start and end of this row
    fStart = fEnd; → start owner face at present cell
    fEnd = ownStartPtr[celli + 1]; → start owner face at next cell

    // Get the accumulated neighbour side
    psii = bPrimePtr[celli];

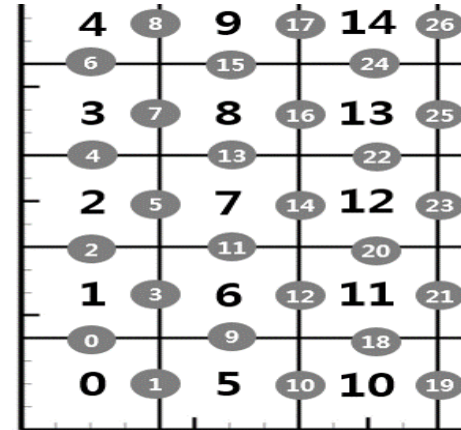
    // Accumulate the owner product side
    for (label facei=fStart; facei<fEnd; facei++)
    {
        psii -= upperPtr[facei]*psiPtr[uPtr[facei]];
    } → psi at cell of owner faces

    // Finish psi for this cell
    psii /= diagPtr[celli];

    // Distribute the neighbour side using psi for this cell
    for (label facei=fStart; facei<fEnd; facei++)
    {
        bPrimePtr[uPtr[facei]] -= lowerPtr[facei]*psii;
    } → psi at cell of neighbour faces

    psiPtr[celli] = psii;
}

```



cellIndex	faceIndex	
	fStart/ownStart	fEnd
0	0	2
1	2	4
2	4	6
3	6	8
4	8	9
5	9	11
6	11	13

Code Analysis of GS

Laplace-Trans

$$F(p) = \int_0^{\infty} e^{-pt} f(t) dt$$

```

scalar psii;
label fStart;
label fEnd = ownStartPtr[0]; → start owner face at each cell

for (label celli=0; celli<nCells; celli++)
{
    // Start and end of this row
    fStart = fEnd; → start owner face at present cell
    fEnd = ownStartPtr[celli + 1]; → start owner face at next cell

    // Get the accumulated neighbour side
    psii = bPrimePtr[celli];

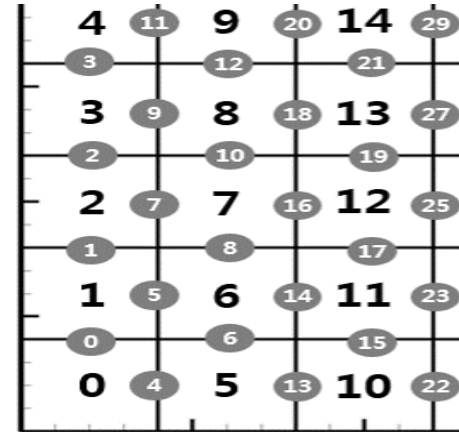
    // Accumulate the owner product side
    for (label facei=fStart; facei<fEnd; facei++)
    {
        psii -= upperPtr[facei]*psiPtr[uPtr[facei]];
    } → psi at cell of owner faces

    // Finish psi for this cell
    psii /= diagPtr[celli];

    // Distribute the neighbour side using psi for this cell
    for (label facei=fStart; facei<fEnd; facei++)
    {
        bPrimePtr[uPtr[facei]] -= lowerPtr[facei]*psii;
    } → psi at cell of neighbour faces

    psiPtr[celli] = psii;
}

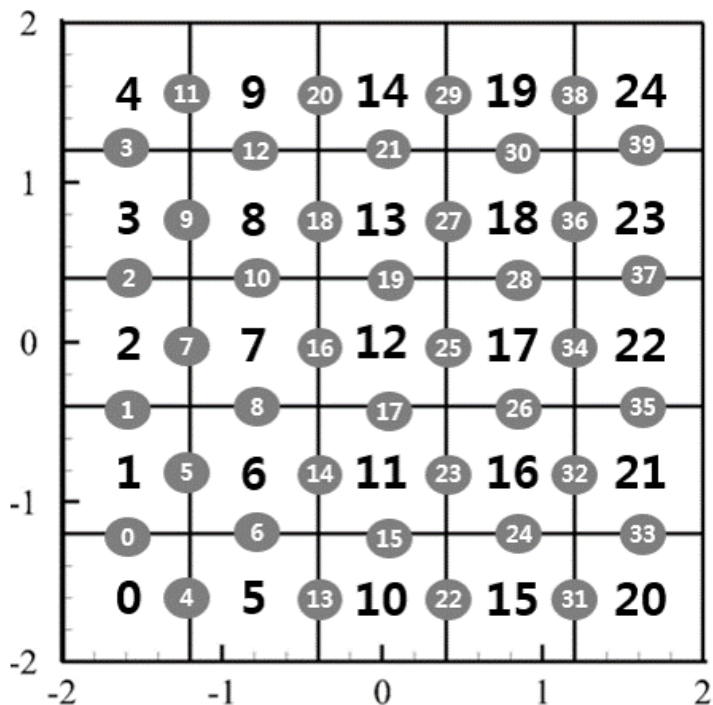
```



cellIndex	faceIndex	
	fStart/ownStart	fEnd
0	0	1
1	1	2
2	2	3
3	3	6
4	6	6
5	6	8
6	8	10

Stabilization of GS

Laplace-Trans
 $F(p) = \int_0^{\infty} e^{-pt} f(t) dt$



```
for(label celli=0;cell<nCells;celli++) psiPtr[celli] = bPrimePtr[celli]/diagPtr[celli];
for(label facei=0;facei<nFaces;facei++)
{
    psiPtr[lPtr[facei]] -= upperPtr[facei]*psiPtr[uPtr[facei]]/diagPtr[lPtr[facei]];
    psiPtr[uPtr[facei]] -= lowerPtr[facei]*psiPtr[lPtr[facei]]/diagPtr[uPtr[facei]];
}
```

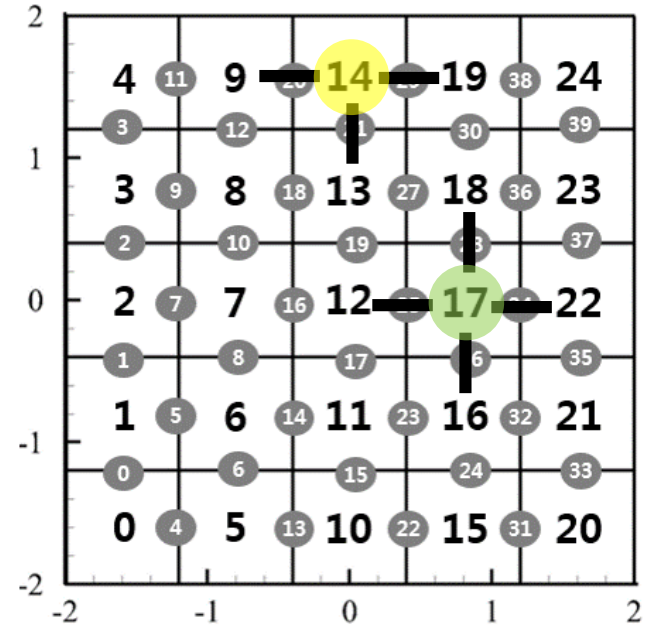
$$\varphi_{i,j} = (l_{i-1}\varphi_{i-1,j} + u_{i+1}\varphi_{i+1,j} + l_{j-1}\varphi_{i,j-1} + u_{j+1}\varphi_{i,j+1})/D_i$$

Stabilization of GS

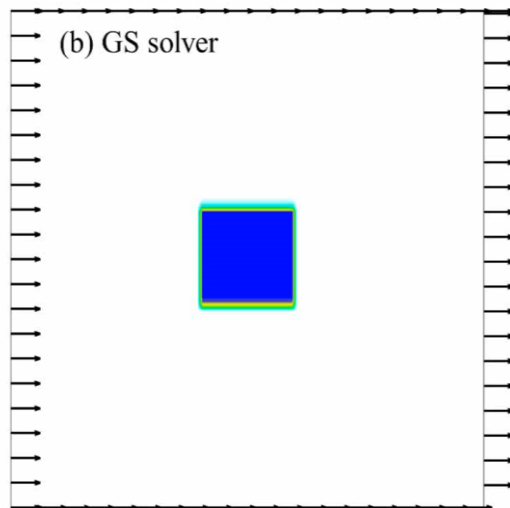
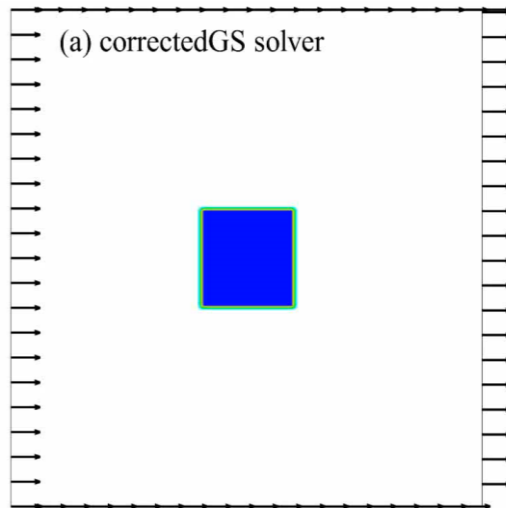
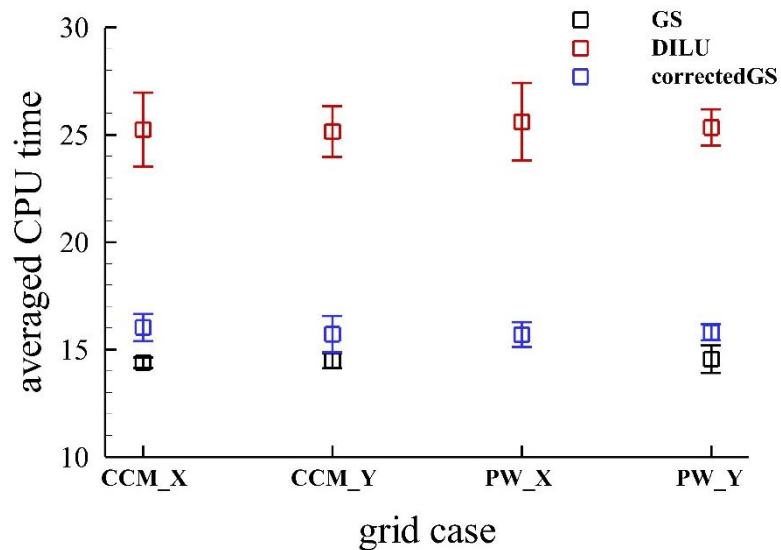
Laplace-Trans

$$F(p) = \int_0^{\infty} e^{-pt} f(t) dt$$

	0	1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	16	17	18	19	20	21	22	23	24	
0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
1	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
2	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
3	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
4	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
5	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
6	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
7	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
8	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
9	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
10	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
11	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0	0
12	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0	0
13	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0	0	0
14	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0	0
15	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.208	0	0	0	0	0	0	0	0	0
16	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.208	0	0	0	0	0	0	0	0
17	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.208	0	0	0	0	0	0	0
18	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.208	0	0	0	0	0	0
19	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.208	0	0	0	0	0
20	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.2	0	0	0	0
21	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.2	0	0
22	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	0	3.2
23	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	0	3.2
24	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	-0.008	0	0	0	0	3.2



Performance of correctedGS



- GS solver
efficient but unstable depending on face-ordering
- DILU solver
stable but inefficient
- correctedGS solver
efficient and stable