

# HPC methods for high resolution Radon transform and deblending

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CREWES Annual Meeting

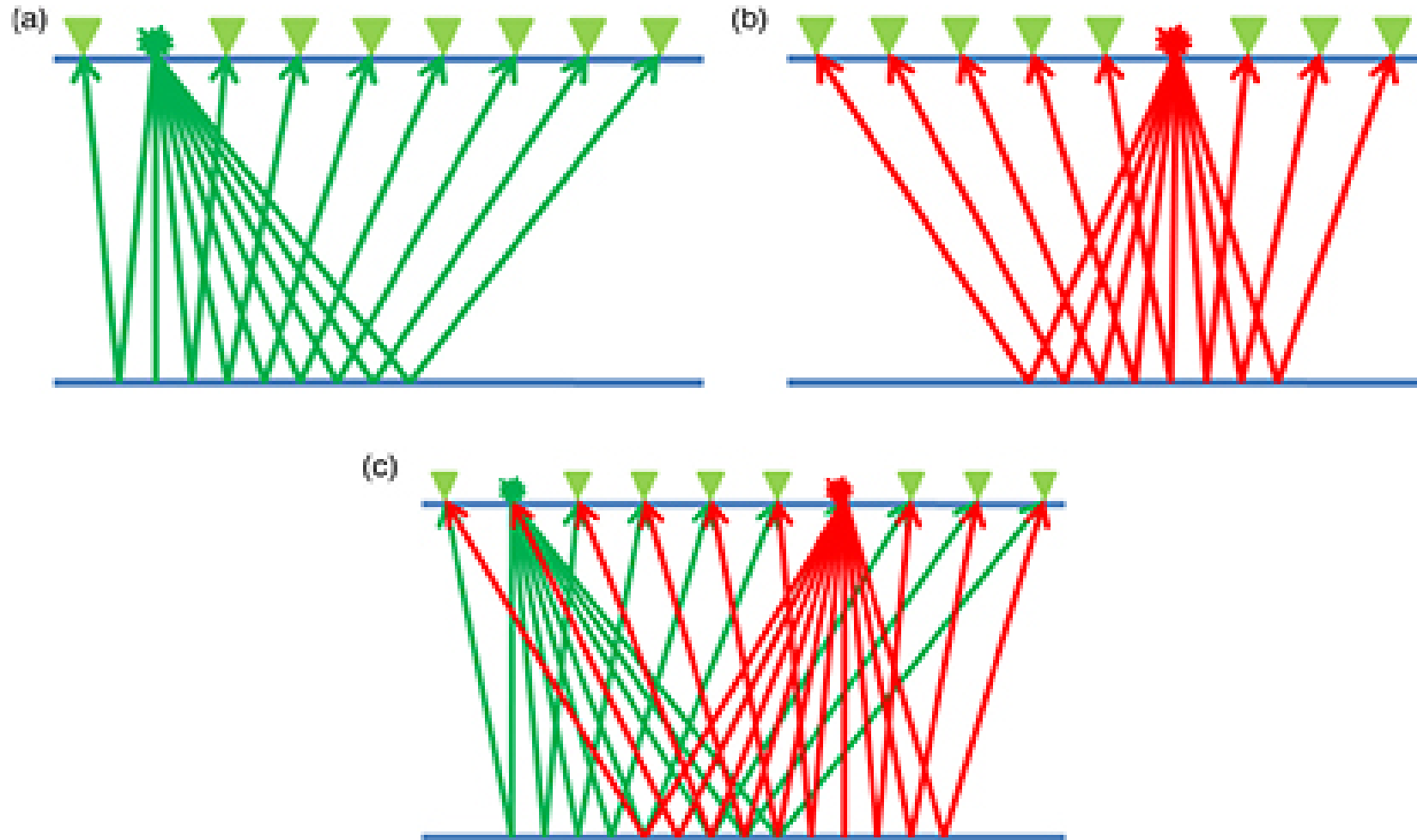
Dec 3 2020



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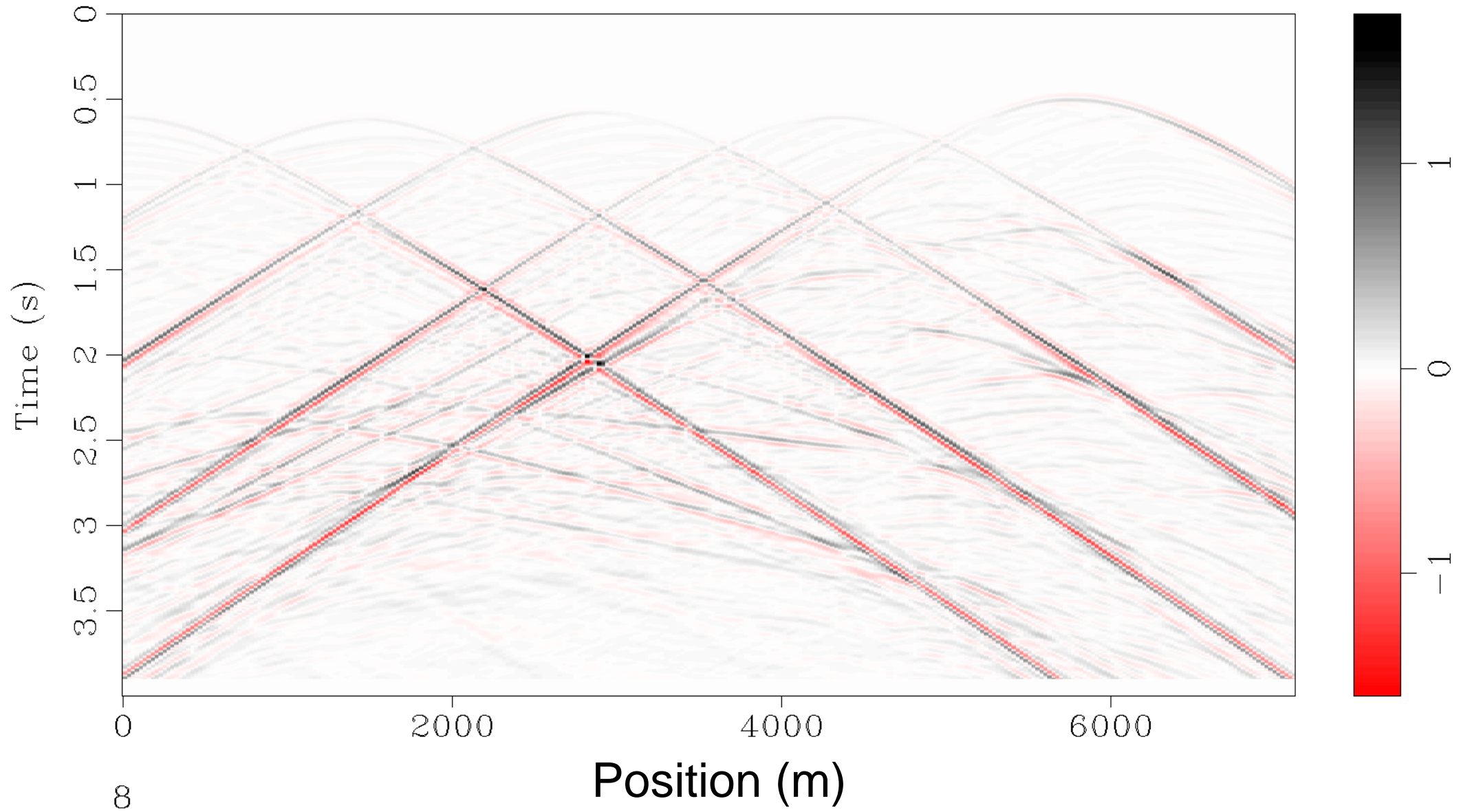


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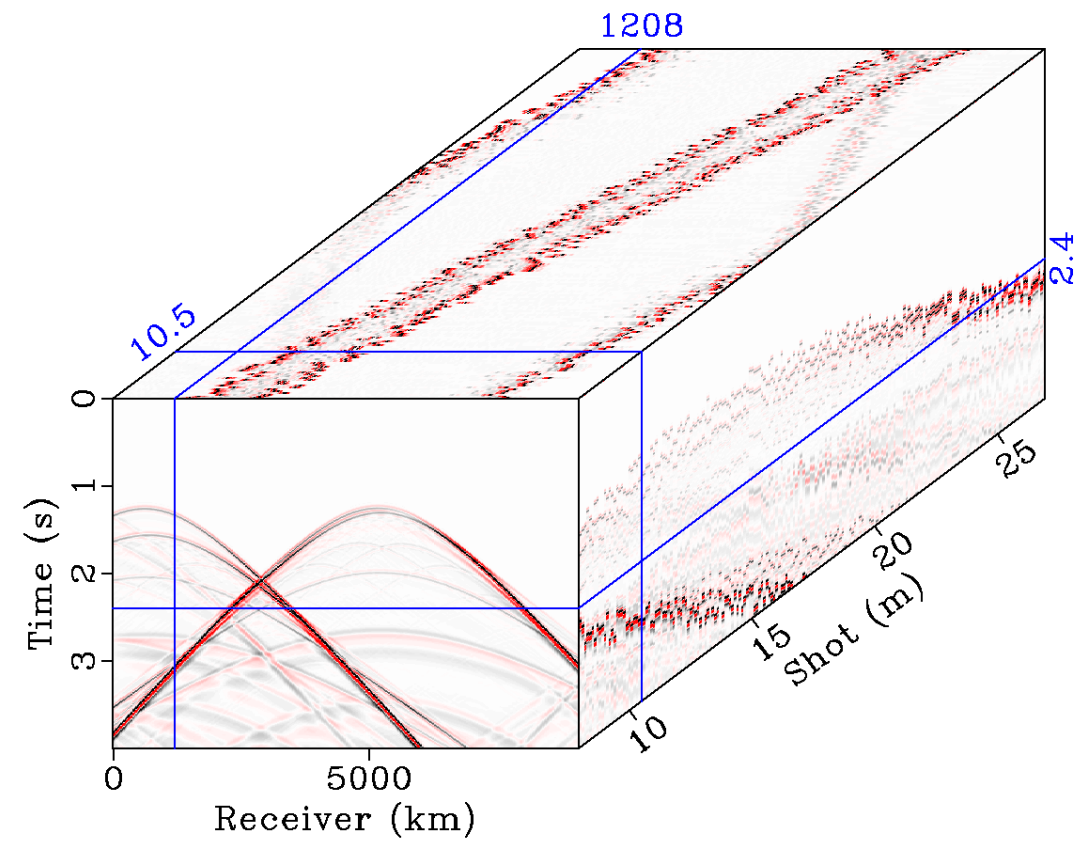


## Blended Marmousi shot

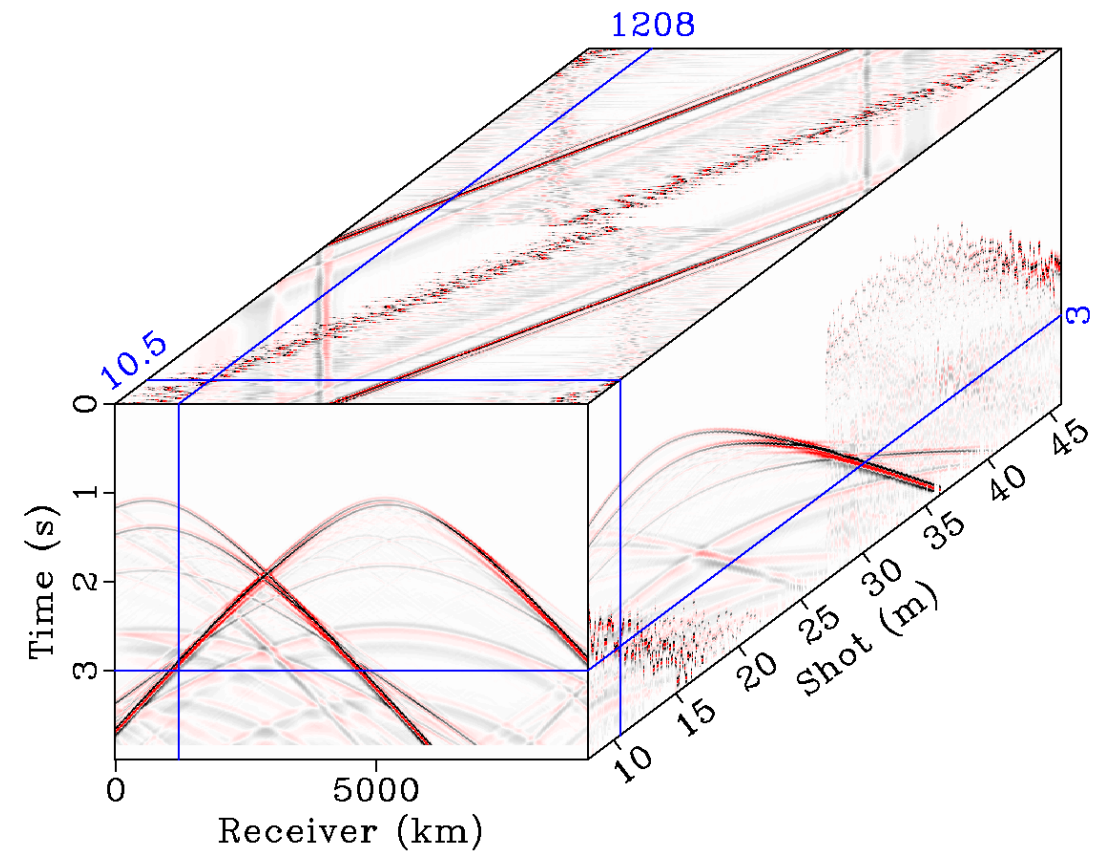




## Blended Data



## Pseudo Deblended Data

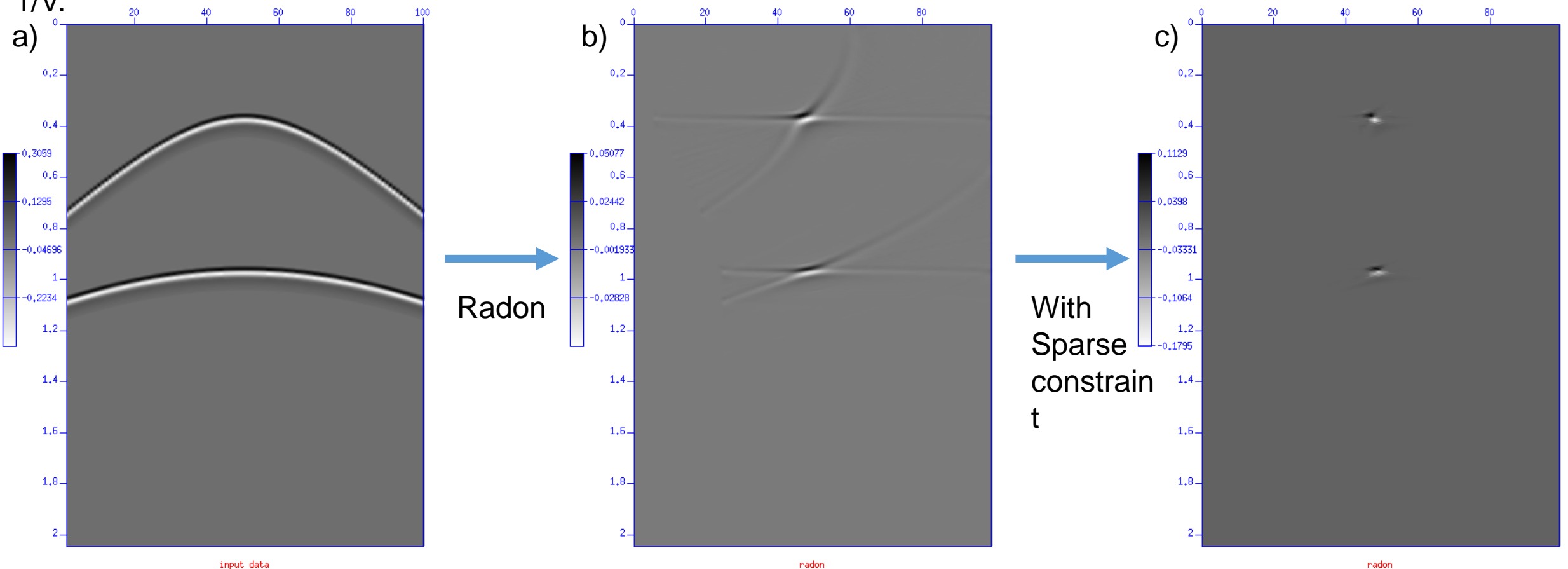




# Sparse Hyperbolic Radon Transform

$$u(p, \tau) = \int_{h_1}^{h_2} d(h, t = \sqrt{\tau^2 + p^2 h^2}) dh$$

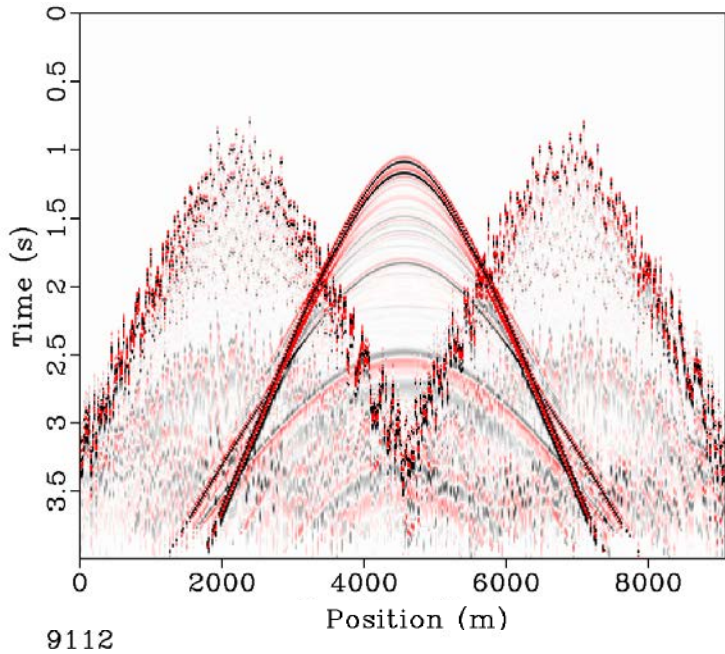
where  $u(p, t)$  is the radon space data,  $p$  is the slowness,  $t$  is the two way travel time,  $h_1$  is the upper offset limit,  $h_2$  the lower offset limit, and  $d$  is the data space to be transformed. The slowness  $p$  is then defined as the inverse of velocity  $1/V$ .



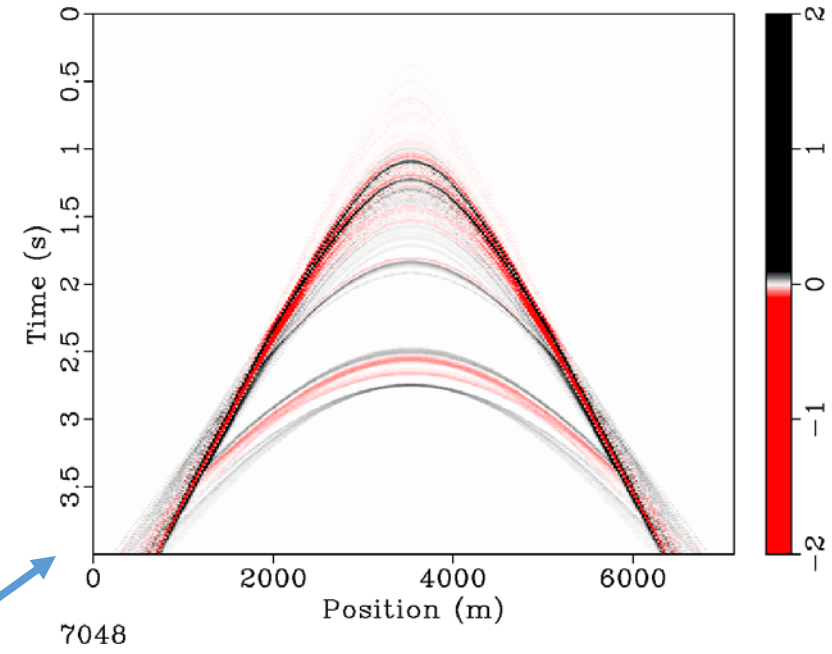
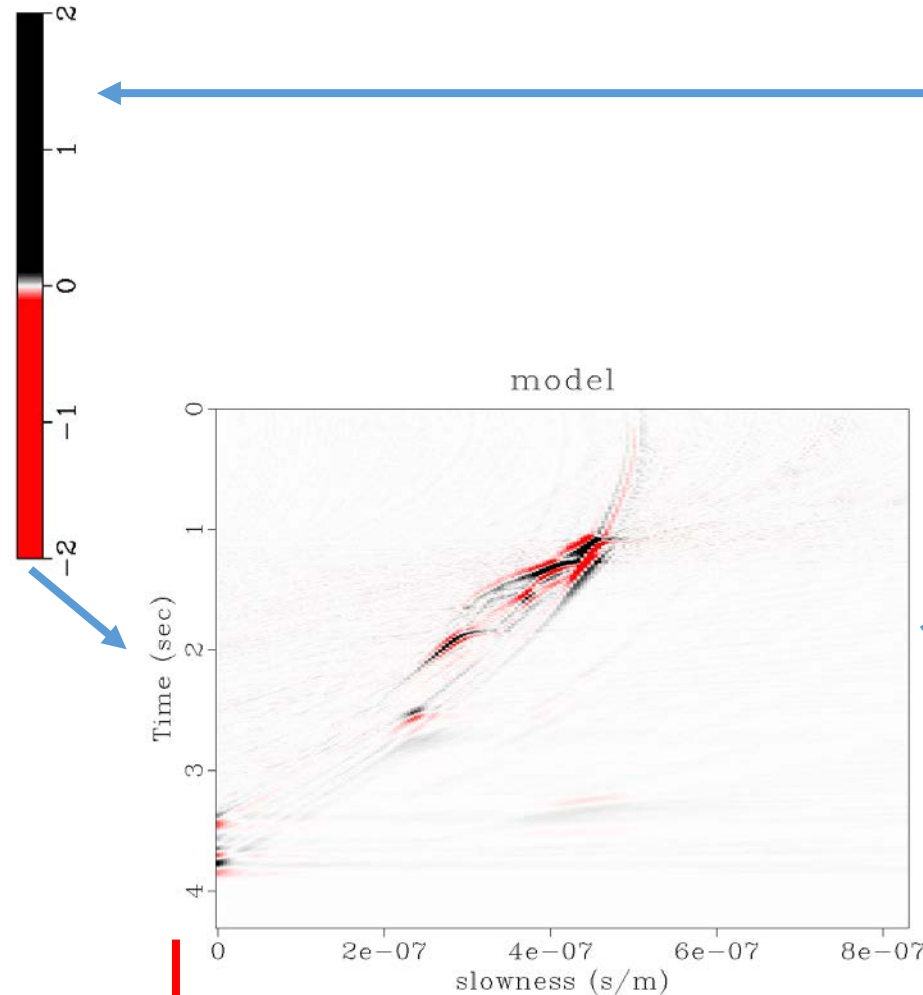


# Denoising – sparse radon transform

$$\left\| S_{pdb} - \boxed{R}m \right\|_2^2 + \mu \|m\|_1$$



Adjoint Operator




Forward Operator



## Radon Denoising

$$S_{pdb} = S_{bl} \Gamma^H$$


$$\left\| S_{pdb} - Rm \right\|_2^2 + \mu \|m\|_1$$

## Radon Inversion

$$\left\| S_{bl} - \Gamma Rm \right\|_2^2 + \mu \|m\|_1$$

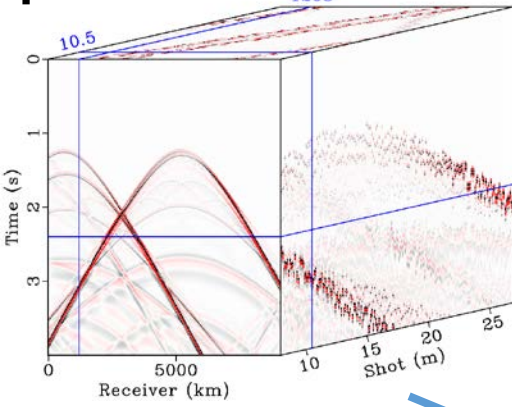




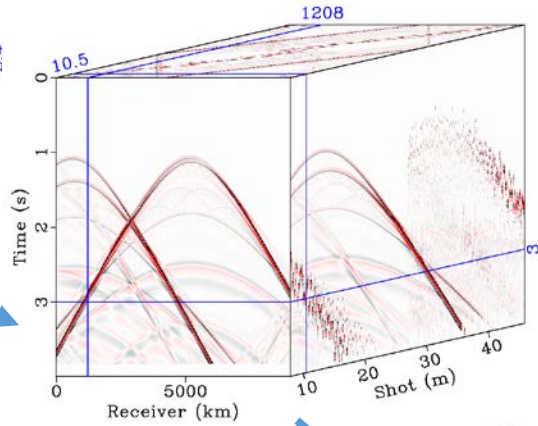
# Sparse Inversion

$$\|S_{bl} - \Gamma R m\|_2^2 + \mu \|m\|_1$$

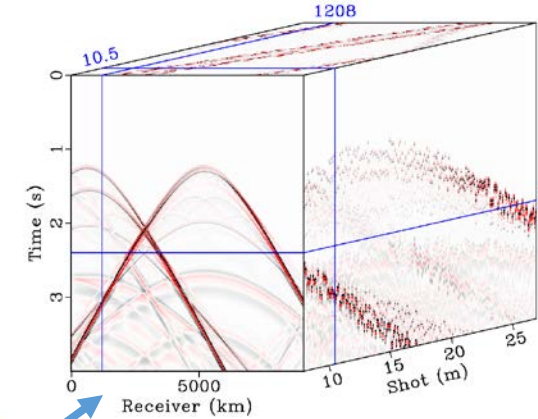
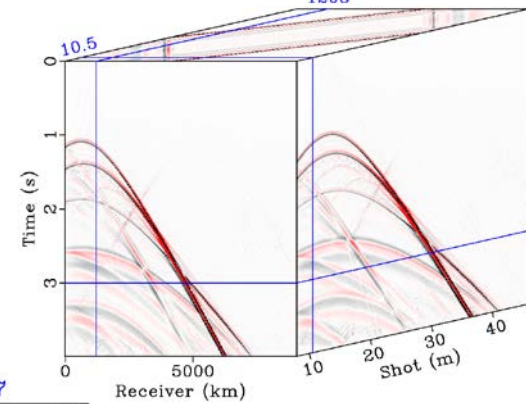
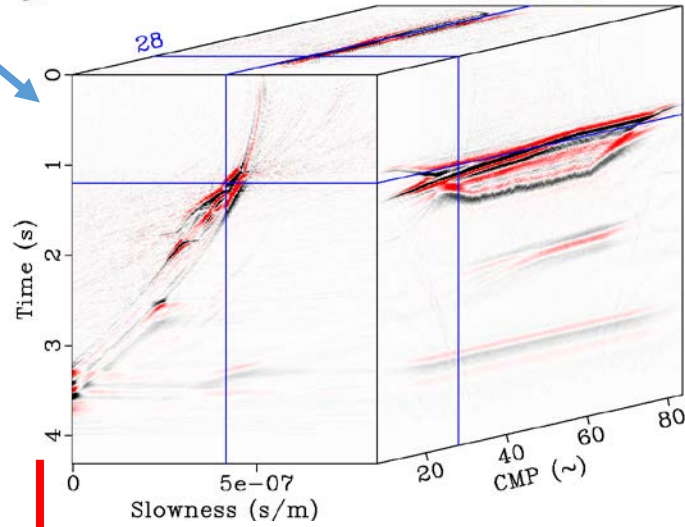
Fitting



Pseudo



Radon



Blending

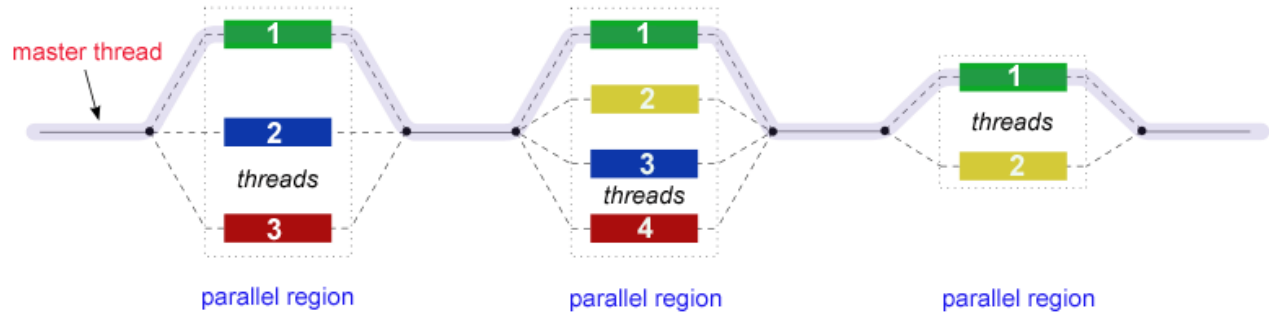
Adjoint Operator

Forward Operator



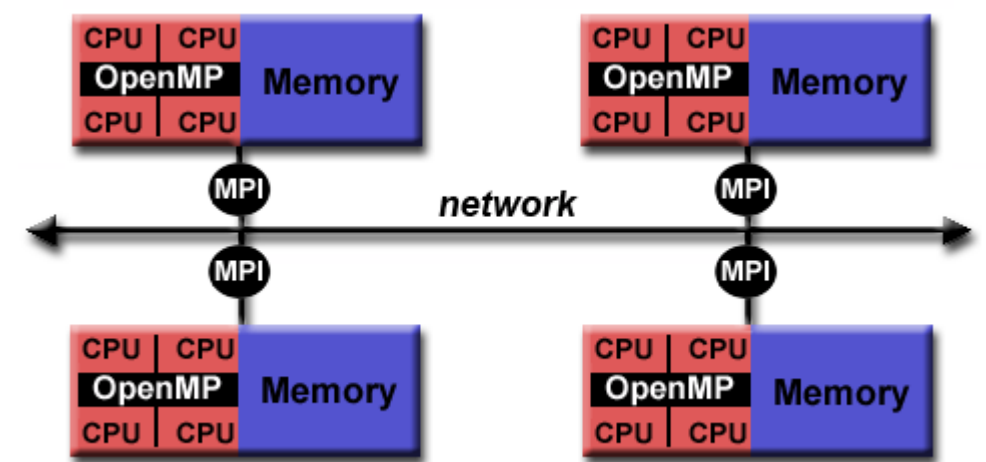
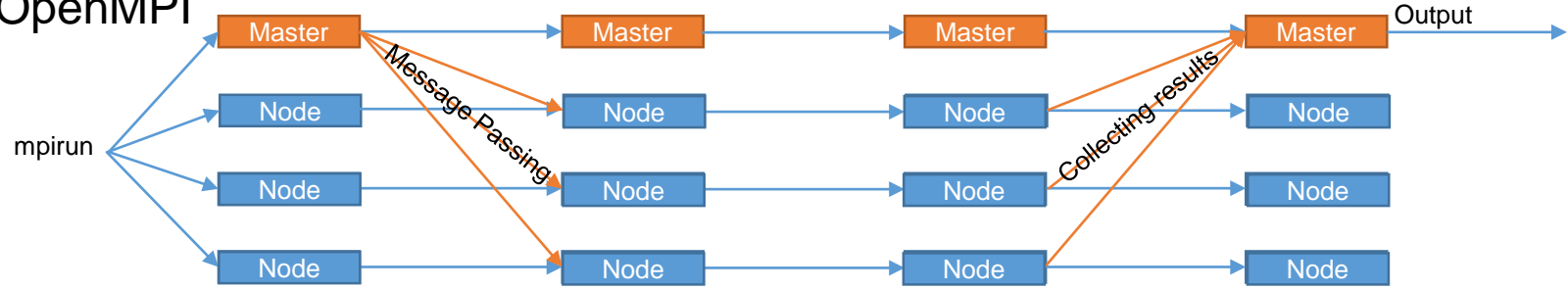
# Parallel programming methods

## OpenMP

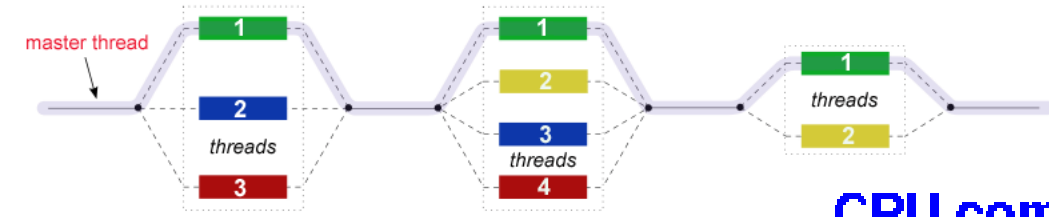


Retrieved from: <https://computing.llnl.gov/tutorials/openMP/>

## OpenMPI

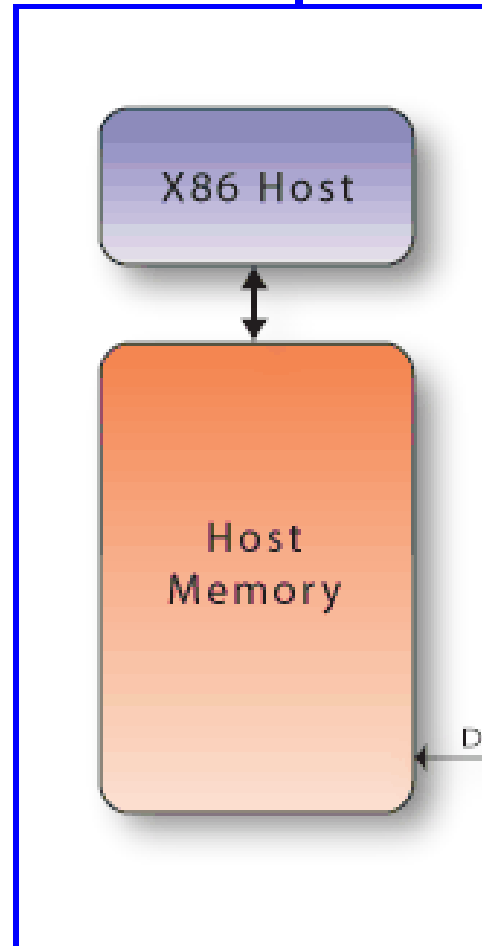


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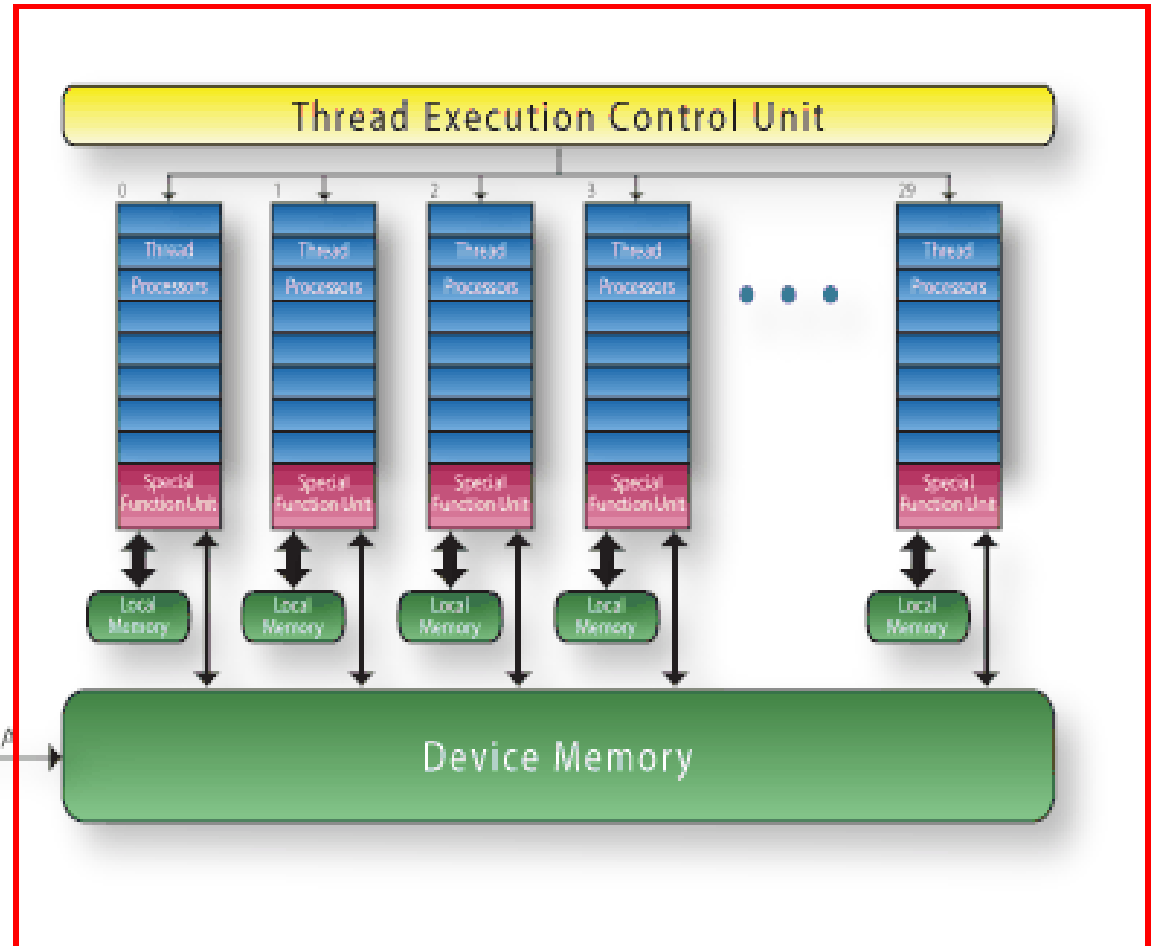
Retrieved from: <https://computing.lnl.gov/tutorials/openMP/>

## CPU computer



## "Traditional" computer

## GPU computer





Key things to note:

## OpenMP

- Simplest to implement limited speedup due to overhead in single threaded portions
- Works on practically any multi-thread CPU

## OpenMPI

- Most complex to implement requiring extensive modification to pre-existing code
- Designed originally for multi-server applications
- Can be used on a single system rather than a cluster for significant performance boost in many scenarios
- Main limitation is excessive data transfers slows execution significantly

## CUDA

- Very easy to implement but requires a Nvidia GPU
- User must be aware of architecture nuances and limitations
- Most effective when memory access is sequential



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**Algorithm 1** Radon pseudo code

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```
1: function RADON
2:   #PRAGMA OMP PARALLEL FOR           ▷ insert for parallelization of loop
3:   for q = slowness do
4:     for h = offset do
5:        $moveout = h^2 * q$ 
6:       for it = 0 to nt do
7:          $time = \sqrt{(it * dt)^2 + moveout}$ 
8:          $model[iqNt + it] += data[ihNt + INT(time/dt)]$ 
9:       end for
10:    end for
11:  end for
12: end function
```

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**Algorithm 2** CUDA Radon pseudo code

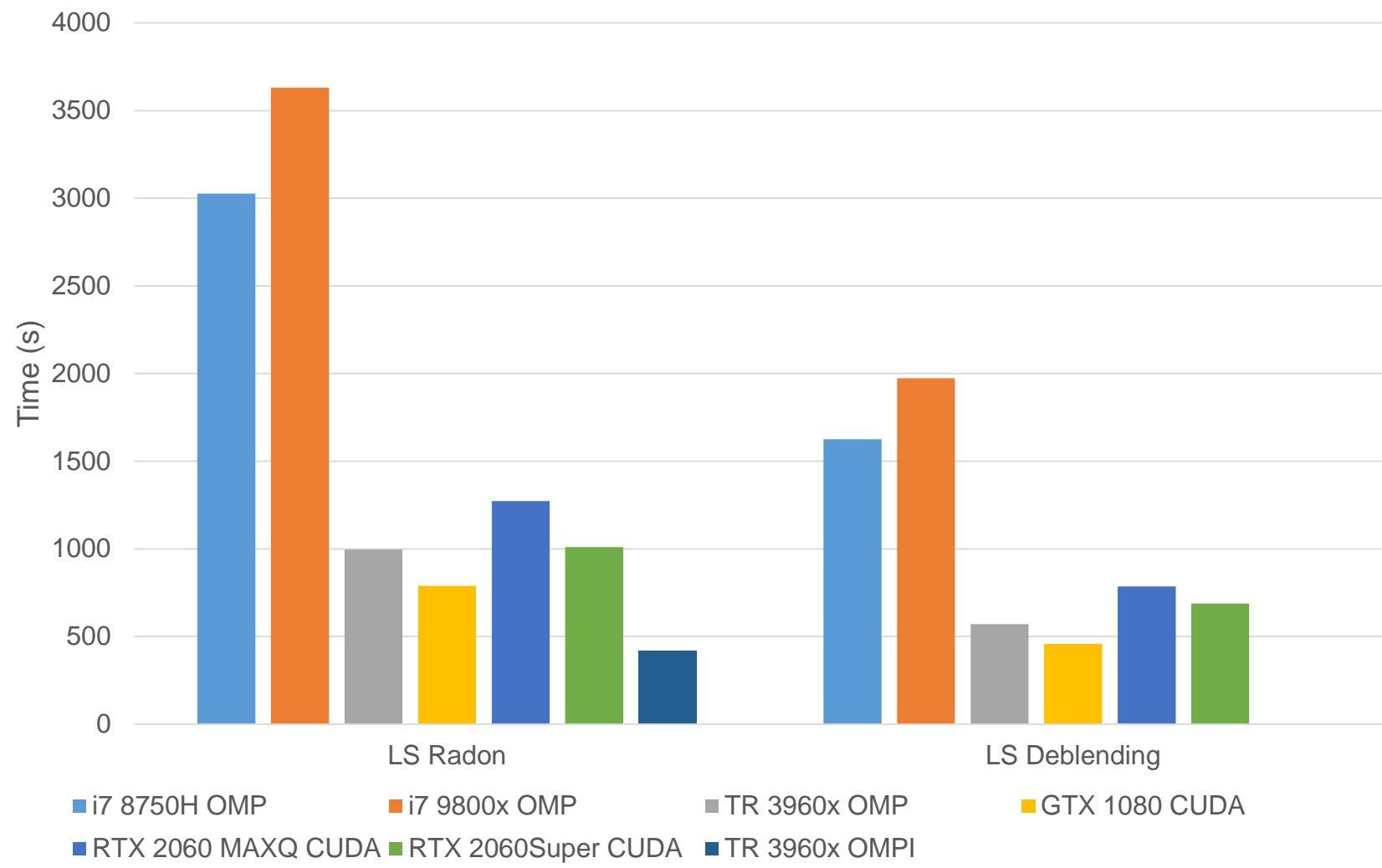
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```
1: __global__
2: function CUDA_RADON
3:   int it = blockIdx.x * blockDim.x + threadIdx.x
4:   int iq = blockIdx.y * blockDim.y + threadIdx.y
5:   int ih = blockIdx.z * blockDim.z + threadIdx.z
6:   if (iq >= nq || ih >= nh || it >= nt) return
7:   int ihNt = ih*nt
8:   int iqNt = iq*nt
9:   double timemax = dt*nt
10:  double  $moveout = h^2 * q$ 
11:  double  $time = \sqrt{(it * dt)^2 + moveout}$ 
12:   $atomicAdd(\&model[iqNt + it], data[ihNt + INT(time/dt)])$ 
13: end function
```

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## Comparative processing time





# Conclusion

- All APIs have good points and draw backs
- CUDA is easy to implement with significant computational advantages for scientific processing
- GPUs are easier to upgrade and price to performance is better than CPU
- Multi GPU is much easier than multi CPU
- openMP is still easiest to implement and is recommended for small workloads
- openMPI is useable for single system parallelization but uses significant amounts of RAM due to allocation per node





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