

Introduction to Computer Science Lecture 04

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Sorting algorithms

- Iterative or recursive,
- Comparison-based or non-comparison based,
- Stable or not stable,
- In-place or requiring more memory

Why review sorting algorithms?

- Usefulness - it's much easier to work on sorted data.
- Sorting algorithms implement many approaches to solve problem.
- Lots of applications:
 - Detecting duplicates,
 - Counting frequency of symbols,
 - Finding subsets,
 - ...and colliding/joining them,
 - Faster searching

Thing we already know...

- Can we use a BST to sort items?
 1. Insert items to BST
 2. Traverse the BST in order.
- PROBLEMS:
 - If a tree is not balanced, and we have a descending order of adding items, we will get a singly linked list instead $\rightarrow O(n^2)$.
 - If we use self-balancing trees, we may enhance this to $O(n \log(n))$.
 - We need $O(n)$ of memory space for it.

Bubble sort

- We have a n -element unsorted/partially sorted array.
- The procedure:

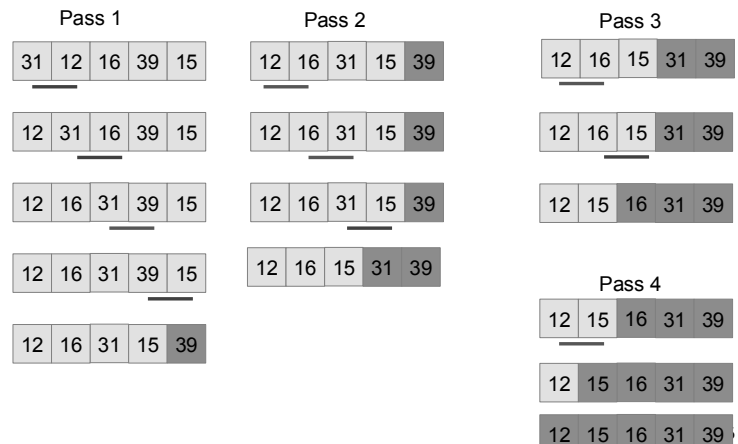
For $x [0..n-1]$

1. Compare pair of numbers n and $n+1$
2. If out of order \rightarrow swap them.

3. Reduce $n-1$ by 1 and repeat.

- Notice how the biggest items „bubble” to the last positions of the array.

Bubble sort



Bubble sort

- Can be implemented using two nested loops.
 - The outer loop spins exactly n times.
 - The inner loop's iterations decrease by 1 with each iteration of the outer loop.
- The complexity is $O(n^2)$

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Let's cheat a little

- Notice that the last pass was done without any swapping.
- If the set is sorted earlier, we will just fruitlessly and blindly compare the already sorted array.
- A good way to make things short is to **terminate the algorithm** when we see that the operation is completed.
- The test will look like this:
 - If during the inner loop iteration no swapping has been made → end the algorithm as the set is sorted.
- This will save some time.

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An altered version complexity

- Worst case is the array sorted backwards, our condition will not run and it will be still $O(n^2)$.
- Best case: A fully sorted array. Will end in a single pass - $O(n)$.
- Practical application: We have to add the (small) performance impact of additional variable holding information was there any swapping or not.

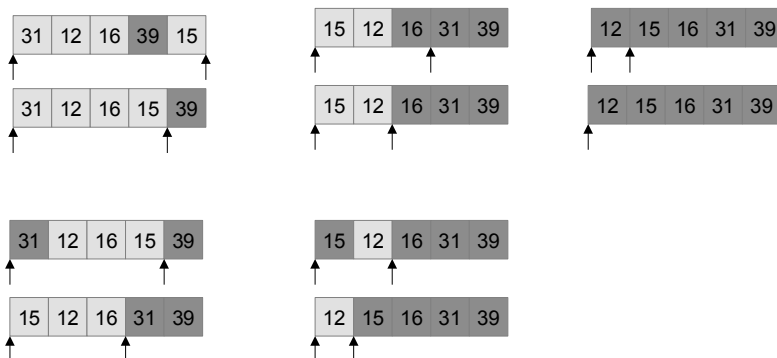
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Selection Sort

- Find the largest item in $0..n$ -element array.
- Swap the largest item with the n -th item
- It's in a proper place, so $n=n-1$ and repeat.

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Selection sort



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Selection sort

- The most important part is finding the index of the last item from the sub-array $[0..n-x]$
- The x is the iteration number then.
- So we still use two nested loops:
 - Outer: Thru a whole array.
 - Inner - Thru less and less elements in the array.
- So the complexity is still $O(n^2)$.

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Implementation considerations

- If our „swap“ is more resource-consuming (we're not using pointers, we have to re-calculate something) the selection sort will significantly outperform bubble sort.
- ...but we cannot use the „cheating“ we used with bubble sort, so we won't get $O(n)$ in the best case.
- In practical applications, usually selection sort performs a bit better.

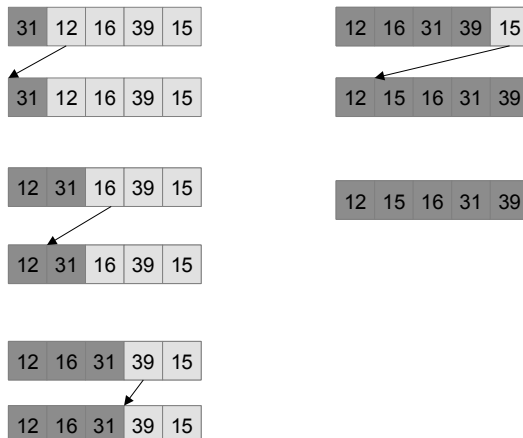
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Insertion sort

- Start with a single item. It is always sorted.
- Take the next item
- Determine the position in the „sorted“ set in which it has to be inserted into.
- Insert the item in its proper position in the „sorted“ set.
- Repeat for all unsorted items.

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Insertion sort



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Insertion sort: Implementation

For every item of index i in the n -item array:

insertedItem = array[i] ← the item to be inserted.

$j = i - 1$

while ($j > 0$) and (array[j] > insertedItem)

array[$j + 1$] = array[j] ← shift sorted items to make space for a new one.

$j--$

array[$j + 1$] = insertedItem ← insert the item in the correct location.

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Insertion sort: Complexity

- The outer loop always executes $n - 1$ times.
- If the array is already sorted, the inner loop will execute once per item.
- In the worst case, insertion will always occur - the loop will execute always at its full range.
- So the best-case complexity is $O(n)$, and the worst-case $O(n^2)$.

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Sorting algorithms stability

- The sorting algorithm is stable when it does not change the relative order of two items with the same value.
- So, with bubble sort, if we swap only if items are in the wrong order, we get the **stable** algorithm. The same with insertion sort.
- Now, the selection sort is **not stable** - starting the largest item lookup from the beginning and inserting it to the end of the sorted set, it will **swap** the order of the largest items with the same value.

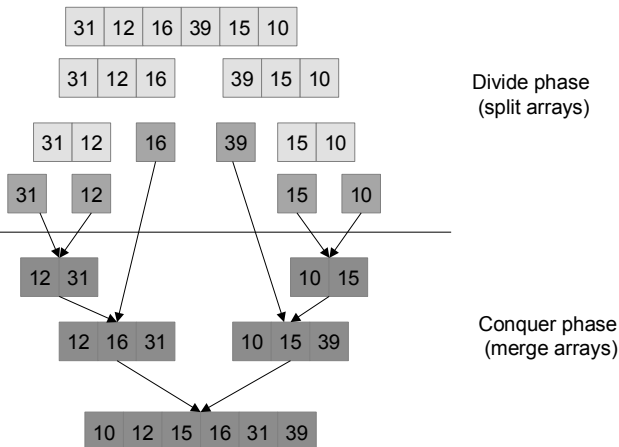
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- Assume we know how to **merge** two sorted sets into one sorted set:
 - $\{2, 3, 10\}$ and $\{5, 17\} \rightarrow \{2, 3, 5, 10, 17\}$
- Can we use it to sort any set?
- We can divide sets as we want.
- A set of 1 item is always in order.
- Now we can **merge 2** 1-item sets.
 - Then we **merge 2** 2-item sets,
 - Then we merge 2 4-item sets,
 - Then we merge 2 8-item sets,
 - » ...
- Until we get a single sorted set.

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- First, we divide the problem to the smaller ones.
- Recursively solve the smaller problems.
- Combine the results of the solutions while coming back from the recursion to obtain solution for larger problem.
- So in the Merge sort:
 - Divide the set to two (equal) halves.
 - Recursively Merge sort two halves.
 - Merge two halves of the sorted array.

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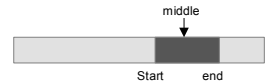
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```

MergeSort(array, start, end)
  if (start < end)
    middle = start + end / 2
    MergeSort(array, start, middle)
    MergeSort(array, middle + 1, end)
  Merge(array, start, middle, end)

```

Is there 0 or 1 item?



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```

Merge(array, start, middle, end)
  n = end - start + 1
  buffer = n-element array
  left = start
  right = middle + 1
  index = 0

  while (left <= middle and right <= end)
    if (array[left] <= array[right])
      buffer[index++] = array[left++]
    else
      buffer[index++] = array[right++]

```

This part inserts new items to the buffer, item by item, it chooses left-hand or right-hand side to insert from.

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```

while (left <= middle)
  buffer[index++] = array[left++]
while (right <= end)
  buffer[index++] = array[right++]

```

All which remains in source sets is copied to the buffer.

```

for every element with index i in buffer
  array[start+i] = buffer[i]

```

Finally, the buffer is copied to the respective place in the array

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- Most of work is done in this Merge function.
- For a call of Merge(array, start, middle, end) we have:
 - start-end+1 items
 - At most *number of items*-1 comparisons
 - At most *Number of items* moves to the buffer
 - ...and the same number or moves back to the main array.
 - So at most it's $3 * \text{number of items} - 1 \rightarrow O(n)$.
- But we call it **many** times...

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- For the single array of n items, we don't call the function.
- When it's divided to 2 sets, we have a single call with $n/2$ items in each half.
- When it's divided to 4 sets, we have 2 calls, $n/2^2$ items in each half (at most).
- When it's divided to 8 sets, 2^2 calls, $n/2^3$ items in each half.
- The total time complexity of the method is $O(n \log(n))$.

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- This is considered an optimal comparison-based method.
- Can operate on large data, requiring the buffer.
- It can be proven, that it is stable.
- ...but it is quite problematic to implement at first (recursion!).
- ...and is not an **in-place** method → requires a non-constant size of extra storage.

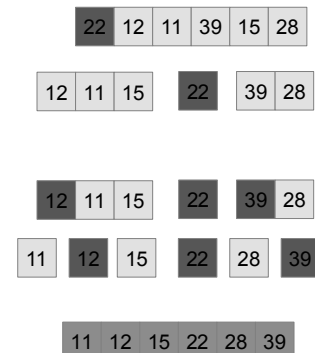
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- In the merge sort, we do most activities in the merge step.
- So in the „divide and conquer“ methodology, we do most of activities after the problem has been divided to sub-problems.
- Is it possible to shift the sorting activity into the divide part?

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- Dividing:
 - Choose a specific item **p** known as **pivot**, and divide the set to two subsets:
 - Smaller than **p**,
 - Larger or equal than **p**.
 - Perform the same thing for every part
- ...and nothing left to do after the recursive division.

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- Notice that the pivot, after executing the sorting round related to it, takes its final position in the given sub-set → it does not participate in further sorting and is in the proper position in this sub-set.
- So the implementation will be:

```
QuickSort(array, start, end)
  if (start < end)
    int pivotIndex = partition(array, start, end)
    QuickSort(array, start, pivotIndex-1)
    QuickSort(array, pivotIndex+1, end)
```

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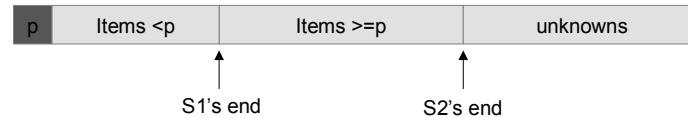
- Assume we will take the 0th item of the array to be partitioned.
- We define two sets (dynamic arrays for example)
 - S1 - in which items are < pivot's value.
 - S2 - in which items are >= pivot's value.
- Iteratively we compare all items against the pivot. If they are smaller, they got to S1, else → S2.
- Finally, we return the S1, pivot and S2.

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- Instead of allocating the memory for two sets, we can **use the array we have**.
- The pivot is array[0] as usual.
- The next parts of the array can be:
 - The first set (S1): < pivot.
 - The second set (S2): >= pivot
 - The last set: Not checked yet.
- We iterate thru items starting from array[0].

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- We iterate on the array starting from pivot.
- If the item is >=p, we increment the pointer to the **S2 set's end**.
- Else, we have to increment the S1's end pointer, but we have an item in ther S2's range...
 - ...so let's just swap these items and extend S2's end pointer too.



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```
Partition(array, start, end)
  int pivot = array[start]
  int S1end=start
  int S2end=start+1;

  while (S2end<=end)
    if (array[S2end]<pivot) //extend S1
      S1end++
      swap(S2end, S1end)
    S2end++ //we always proceed with S2

  swap(start, S1end) //correct pivot position
  return S1end //return new pivot position
```

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- The algorithm is **not stable**, but **in-place**.
- The best result is obtained if the problem is always divided to two equal halves.
 - Then, depth of recursion is logarithmic,
 - The complexity is $O(n \log(n))$.
- The worst case is when the pivot gets always separated → we get a full S1 and empty S2 or full S1 and empty S1.
 - Then, we get to $O(n^2)$
 - Notice this will happen is we try to sort the already sorted array!
 - It can be fixed with different pivot initialization.

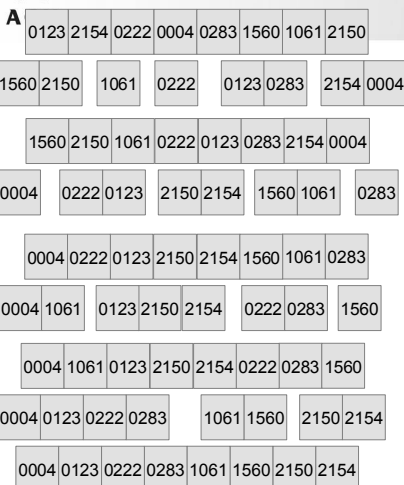
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- We have always used sorting algorithms for numbers.
- What if we want to sort strings?
 - Convert strings to ASCII numerals?
 - ...?UTF-8 numerals? ← trouble!
- What if we want to develop a sorting method **specifically** for strings?

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- We consider each record of data as a string of symbols.
- We group string into sets according to the next symbol in each string.
- Concatenate the sets for the next iteration
- Repeat until sorted.
- Assume we have a constant-length strings.
 - ...but if we stick to numerals, we can zero-pad.

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Grouped by the 4th symbol....

...and concatenated back.

Grouped by the 3rd symbol...

...and concatenated back.

Grouped by 2nd symbol...

...and concatenated back.

Grouped by the 1st symbol...

...and it's a sorted set. 39



- How do we group items?
 - We find the first item of the specified characteristics,
 - Move it to the specific set.
 - Go until the end of input data.
 - Repeat for each symbol (or group).

So:

`i=1;`

for every symbol in the record length:

group the items by i-th item.

concatenate groups

`i++`

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A well aligned example from Halim S. - "World of Seven" - <https://www.comp.nus.edu.sg/~stevenha/>



- We can do grouping iteratively:

Given `i` = the position we use for grouping.
Create a set of vectors for every symbol.

for every symbol in the array:
add the symbol to the vector related to it.

- For the decimal numbers:

```
digit=0;
for every element of the array
  digit=(element/i) %10
  push(digitsList[digit], element)
```

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- We can just copy the vectors to the array, symbol by symbol:

`i=0`

for every symbol in the alphabet

while symbol's list not empty

array[i]=pop(symbol's list)

`i++`

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- We can use any alphabet we want.
- For each iteration we go thru the whole array once to place them to groups, then we concatenate groups to the array.
- So the complexity is $O(n)$.
- Number of iterations: number of symbols in the alphabet.
- So the complexity is $O(dn)$.
- Not in-place, but stable.
- Requires more memory.

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- A binary **heap** is a data structure which is a binary tree with the following limitations:
 - All its levels except the last one are completely filled (Shape property).
 - The value stored in each node is greater (or less - depending on implementation) than its children.
- Thanks to shape property, we can **serialize** a BST into the linear array and retrieve it back without any pointers.
- Because it's not a BST, we can shift elements down without any side effects.

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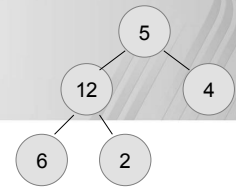
- If we are able to build a heap of the data, we can do the following:

1. Remove the topmost part of the heap (maximum) and append it to the sorted list part.
2. Re-create the heap.

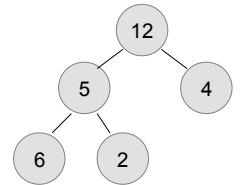
- Because the n-element heap is serialized in the array, we can do the step 1 by moving the maximum to the place right after the end of the heap in the same array - the „heap“ part will shrink, the „sorted“ part will grow then.

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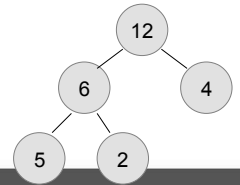
- [5,12,4,6,2]
 - Not a heap



- As $5 < 12$, we swap them



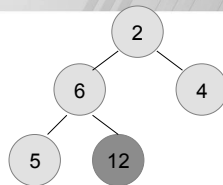
- And as $5 < 6$, we swap 5 with 6.



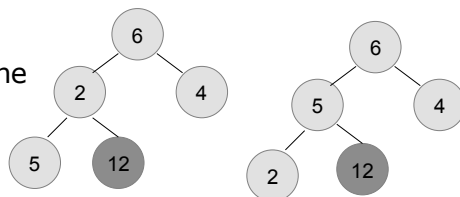
- We have a heap property.

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- Remove the root - this is sorted
- We do it by swapping it with the last node and trimming it off from further processing.

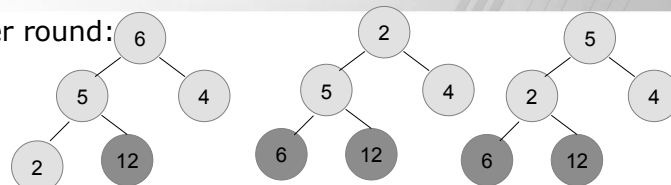


- Then, we re-create the heap:
 - Swap 6-2
 - Swap 5-6



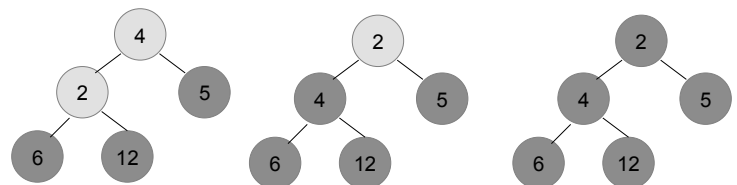
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- Another round:



Pop the element to the end

Heapify



Pop the element

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Heap sort: Typical Implementation

- HeapSort(array, count)


```

start=count/2
end=count
while (end>1)
  if (start>0)
    start--
  else
    end--
    swap(array[end],array[0])

//Moving the smaller elements down
root=start
while (2*root+1)<end
  leaf=2*root+1
  if (leaf+1<end and a[leaf]<a[leaf+1])
    leaf++
  if (a[root]<a[leaf])
    swap(a[root],a[leaf])
    root=leaf
  else
    break
      
```

}

Select the largest item and put it to the end

Restore heap properties.

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Heap sort

- Complexity (worst case): $O(n \log(n))$
- Not a stable sort
- In-place algorithm
- In many practical cases a bit slower than Quick sort.

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Hybrid algorithms

- How can we get rid of $O(n^2)$ complexity in badly formed input data?
 - Use a different algorithm, not so suitable for other alignments of data...
- Introspective sort variant:
 - If the number of data is relatively small, go with InsertionSort.
 - If the recursion depth of Quick sort is acceptable we can go with Quick sort.
 - ...but we can just partition the array in 2 and go Intro Sort on them.
 - Else, we go with Heap sort.

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Estimating the depth of recursion

- Usually $2 * \log_2 n$.
- This will „lock“ the complexity to $O(n \log(n))$.
- The total complexity of Intro sort will be then $O(n \log(n))$
- ...however, it is not stable.
- ...and more difficult to implement.

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A small interruption

- Stack machines
 - A CPU that does not use registers, but (usually one) stack.
 - Execution of operation means:
 - Fetching operation code from the stack,
 - Fetching operands from the stack,
 - Pushing the result to the stack.
 - With 2-operand command, stack is shorter by 1 value or 2 items: instruction and value.

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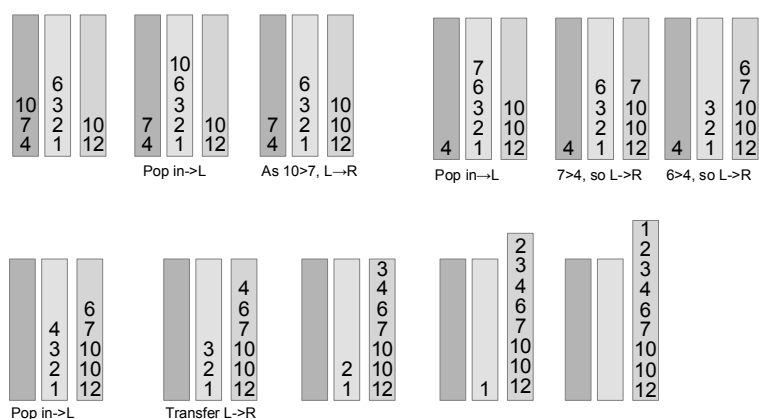
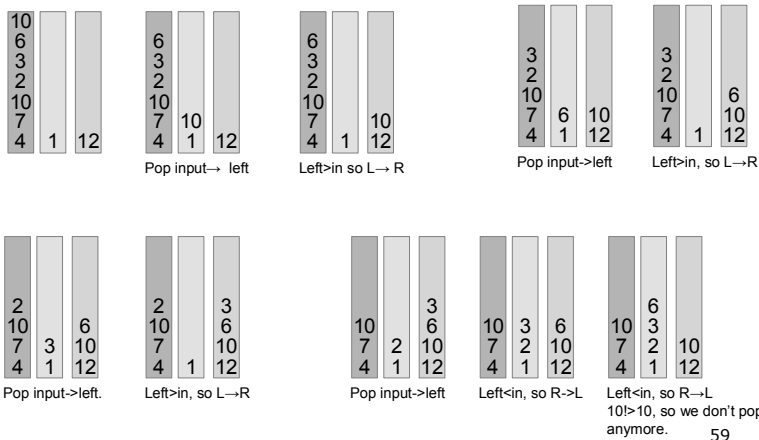
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- Implementations:
 - Separate stack and conventional instructions storage (Symbolics, 1970s).
 - Single stack including instructions and operands (Ferranti, 1965-70s).
 - Multiple stacks, switchable and shorable (Syeika, 1970s, Multiklet project).
- Because it is quite difficult to program these CPUs and typical programming languages compilers don't generate such code efficiently, they never got popularity.
 - In 1970s, Symbolics LISP-programmable machines were implemented as stack CPUs.

- Having a set of stacks it is possible to implement a sorting algorithm similar to insertion sort.
- It can be implemented using conventional programming techniques, but also with a stack-based processors.
- The „register-as-stack“ architecture allow to obtain a very high performance in the implementation.

- Rules:
 - No access different than stack-based,
 - No registers, only stacks,
 - The only moving-related operations are push and pop. Popped value must be pushed somewhere.
 - We can compare top values from various stacks without popping.
- Given:
 - Input, unsorted data stack
 - Two working stacks: Left and right.
 - Minimum and maximum element value.

- Initialization:
 - Push the minimum to the left stack,
 - Push the maximum to the right stack.
- Operation:
 - Pop the value from input stack to the left stack.
 - Until top of left stack is not larger than top of input stack:
 - Pop the left stack to the right stack.
 - Until top of the right stack is not smaller than top of input stack:
 - Pop the right stack to the left stack.



Summing up

- This is an insertion sort.
- Insertion sort can be implemented on stack-like data structures.
- With specific stack machine architectures, it is possible to make it even more efficient.

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Mass-solving linear equations

- Any discretized set of differential equations can be described as a system of linear equations.
- The problem is that number of these linear equations may be VERY large.
- Is it possible to solve these equations with a computer?

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Linear equations system

- A set of linear equations like:

$$\begin{cases} Ax_1+Bx_2+Cx_3+Dx_4=U \\ Ex_1+Fx_2+Gx_3+Hx_4=V \\ Ix_1+Jx_2+Kx_3+Lx_4=W \\ Mx_1+Nx_2+Ox_3+Px_4=X \end{cases}$$

- Can be described using matrix equation:

$$\begin{bmatrix} A & B & C & D \\ E & F & G & H \\ I & J & K & L \\ M & N & O & P \end{bmatrix} \cdot \begin{bmatrix} x_1 \\ x_2 \\ x_3 \\ x_4 \end{bmatrix} = \begin{bmatrix} U \\ V \\ W \\ X \end{bmatrix}$$

- Sometimes described as:

$$\mathbf{A} \times \mathbf{X} = \mathbf{B}$$

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Gauss method

- Allows to obtain solution of such equations system.
- Inspired by the method of solving these systems by eliminating unknowns until we get some $x_n=Y$, a single unknown solved.
 - It can be used to obtain the next equation's unknown,
 - And then, using two solutions, another one...
 - Going this way, we can obtain all unknowns.

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Elimination

- If we get a $A \cdot X = B$ matrix equations (in which X and B are vectors), we can write A|B matrix:

$$\left[\begin{array}{cccc|c} a_{1,1} & a_{1,2} & a_{1,3} & \dots & a_{1,n} & b_1 \\ a_{2,1} & a_{2,2} & a_{2,3} & \dots & a_{2,n} & b_2 \\ a_{3,1} & a_{3,2} & a_{3,3} & \dots & a_{3,n} & b_3 \\ \cdot & \cdot & \cdot & \dots & \cdot & \cdot \\ \cdot & \cdot & \cdot & \dots & \cdot & \cdot \\ \cdot & \cdot & \cdot & \dots & \cdot & \cdot \\ a_{n,1} & a_{n,2} & a_{n,3} & \dots & a_{n,n} & b_n \end{array} \right]$$

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Elimination (2)

- Next, we convert all elements under $a_{1,1}$ to 0.
 - Then all elements under $a_{2,2}$,
 - Then, $a_{3,3}$ etc.
- (the first row remains the same)
- So we can use a single solution to expand it to the rest.

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- Elimination of column 2 (so only $a_{1,1}$ remains):
 - For all elements of the row i ($2..n$) we add the next elements of the row 1 multiplied by $-1*(a_{i,1}/a_{1,1})$.
 - Notice that for $a_{2,1}$, we will get:
 - $a_{2,1} - (a_{2,1}/a_{1,1}) * a_{1,1} = a_{2,1} - a_{2,1} = 0$
 - ...and that's what we want!
 - For $a_{2,2}$ we will get: $a_{2,2} - (a_{2,1}/a_{1,1}) * a_{1,2}$
 - For $a_{2,n}$ we will get: $a_{2,n} - (a_{2,1}/a_{1,1}) * a_{1,n}$
 - For b column: $b_2 - (a_{2,1}/a_{1,1}) b_1$

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- We process it until we will get the matrix **triangular**:

$$\begin{bmatrix} a_{1,1} & a_{1,2} & a_{1,3} & \dots & a_{1,n} & b_1 \\ 0 & a'_{2,2} & a'_{2,3} & \dots & a'_{2,n} & b'_2 \\ 0 & 0 & a'_{3,3} & \dots & a'_{3,n} & b'_3 \\ \cdot & \cdot & \cdot & \dots & \cdot & \cdot \\ \cdot & \cdot & \cdot & \dots & \cdot & \cdot \\ 0 & 0 & 0 & \dots & a'_{n,n} & b'_n \end{bmatrix}$$

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- We will then unpack the matrix to something that looks more like an equation system:

$$\begin{bmatrix} a_{1,1} & a_{1,2} & a_{1,3} & \dots & a_{1,n} \\ 0 & a'_{2,2} & a'_{2,3} & \dots & a'_{2,n} \\ 0 & 0 & a'_{3,3} & \dots & a'_{3,n} \\ 0 & 0 & 0 & \dots & a'_{4,n} \\ \cdot & \cdot & \cdot & \dots & \cdot \\ 0 & 0 & 0 & \dots & a'_{n,n} \end{bmatrix} \cdot \begin{bmatrix} x_1 \\ x_2 \\ x_3 \\ x_4 \\ \cdot \\ x_n \end{bmatrix} = \begin{bmatrix} b_1 \\ b'_2 \\ b'_3 \\ b'_4 \\ \cdot \\ b'_n \end{bmatrix}$$

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- Formula for the unknown x_i :

For $i=n-1, n-2, \dots, 1$:

$$x_i = (b'_i - a'_{i,n}x_n - \dots - a'_{i,i+1}x_{i+1}) / a'_{i,i}$$

- So:

$$x_n = b'_n / a'_{n,n}$$

- Knowing that subsequent operations are recursive, it can be implemented recursively.

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Stage 1: Gaussian elimination

```

For i=1,2,...n-1
  For j=1,2,...n
    If AB[i,i]==0
      return 2 //Error! We are going to divide by 0!
    multiplier = AB[j,i]/AB[i,i]
    For k=i+1..n+1
      AB[j,k]=B[j,k]+multiplier*AB[i,k]
  
```

WARNING!
 ==0 is assumed to
 have some tolerance!

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Stage 2: Obtaining unknown values

```

For i=n, n-1, n-2 ... 1
  s=AB[i, n+1]
  For j=n, n-1, n-2 ... i+1
    s=s-AB[i,j]*x[j]
  if AB[i,i]==0
    return 2 //ERROR - we are going to divide by 0
  X[i]=s/AB[i,i]
  
```

- Unknowns are stored in X vector

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Crout's enhancement

- Presence of any 0 in the diagonal of the matrix, or introduction of such 0, will cause the algorithm to divide by 0.
- Because $A+B=B+A$, we can swap columns in the array as we wish.
- Implementation of column swap can be done just using pointer swap.
- Can we swap columns to not get zeros in the diagonal, or better, get them in part we want?

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- We search for the element with biggest absolute value.
- Now, we swap columns: Column with this element with column with the part of the diagonal.
- It can minimize the division by zero errors.
- Instead of pointers, a **lookup table** can be used.

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Even better version

- In practical applications, frequently we have this $A \cdot x = b$ problem with various b values for the same, or similar, A values.
 - Like the same discretized continuous problems, for the same equations, for different points in the medium.
- We can save the eliminated A -values and use them again.
 - ...but there is a b -vector involved - the decomposition must be different.

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$A = LU$

- We describe the converted A matrix as a matrix product of upper and lower triangular matrices:

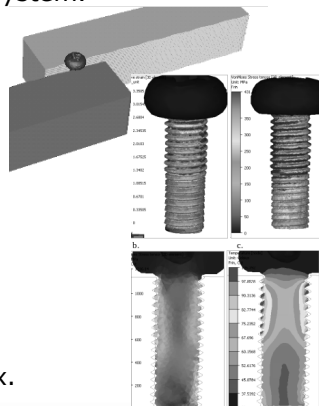
$$A = LU$$

- Once we decompose A to LU , we can save them and substitute with any values of B we want.
- The complexity remains as in Gaussian elimination.

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Why do we need this?

- A lot of physical, mechanical, structural problems can be converted into a linear equations system.
- There are even more efficient methods to do this.
- If we discretize the continuous media and interpolate in between, we can solve non-linear problems using linear equations!
- ...however, the typical method for an e.g. computer simulations has millions of columns and rows.
 - ...fortunately it is a sparse matrix.



Graph algorithms

- A **graph** is a data structure which consists of **vertices** and **edges** linking them.
- Both vertices and edges may hold additional information.
- The **graph order** is a number of vertices in the graph.
- The **graph size** is a number of edges in graph.
- A **null graph** consists vertices, but no edges.
- Graphs are used in discrete mathematics, geometry/topology, and, in application, in engineering.

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Properties of graph elements

- Graph may allow a **multi-edge** connection, it means that two vertices may be connected by more than one relation.
- A **loop** is a connection to itself.
- The edge may be **uni- or bidirectional**.
- The graph is **planar** if we can draw it without crossing the edges.
 - Finding planar graph of an existing graph is an important problem in design software algorithms.

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Properties of graph elements

- Graph's edges may have their values, called **weights**.
- A **path** is a series of vertex traversals from one vertex to another, usually thru other vertices.
 - Finding the shortest path is an important problem in function optimization and logistics.
- A **simple graph** has no multi-edges or loops.

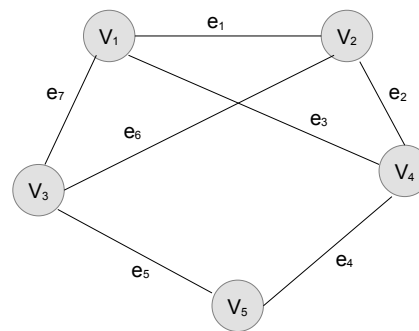
80

Paths and cycles

- A **Hamiltonian path** is a path which goes thru all vertices of the graph.
- A **Hamiltonian cycle** is similarly, a closed path. **The simple cycle must cross each vertex only once.** Some edges may not be used.
- An **Eulerian path** goes thru all edges.
- In the **Eulerian cycle** the path must be closed.

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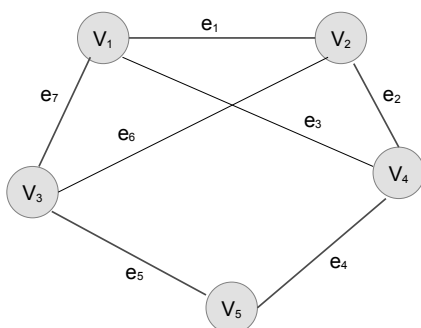
Graph example



A **non-directed, simple but non-planar** graph.
Not a **complete** graph as V_5 is not directly connected to e.g. V_1 or V_2 .

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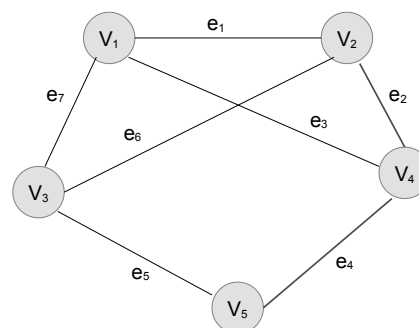
Graph example



A Hamilton cycle

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Graph example



A V_5 to V_2 path.

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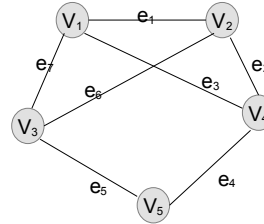
How graph can be stored?

- Structures and pointers
 - PROBLEM: If we may have any number of edges from a vertex, we need dynamic data structure to hold pointers.
 - PROBLEM: Lack of general overview of the graph.
 - PROBLEM: Algorithms which look for specific vertex will have to traverse it almost blindly.

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Better way to store graphs?

- Adjacency matrix:
 - For n vertices we create $n \times n$ matrix of binary values.
 - 1 if there is a connection between column- and row-related element.
 - We can store **directed** and non-directed graphs.

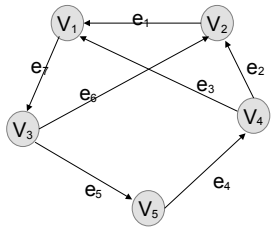


	V ₁	V ₂	V ₃	V ₄	V ₅
V ₁	0	1	1	1	0
V ₂	1	0	1	1	0
V ₃	1	1	0	0	1
V ₄	1	1	0	0	1
V ₅	0	0	1	1	0

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Better way to store graphs?

- Adjacency matrix and directed graphs:
 - Rows contain starting points
 - Columns: End of edge.



	V ₁	V ₂	V ₃	V ₄	V ₅
V ₁	0	0	1	0	0
V ₂	1	0	0	0	0
V ₃	0	1	0	0	1
V ₄	1	1	0	0	0
V ₅	0	0	0	1	0

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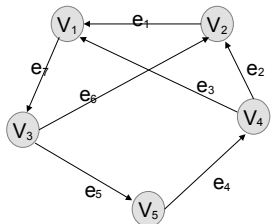
Using the adjacency matrix

- A degree of the vertex, for a non-directed graph, is a count of 1s in the column or row related to this vertex.
- For a directed graph, we obtain, by counting in columns or rows, degree of „inputs“ or „outputs“ of the vertex.
- In a directed graph, if $[x,y] = [y,x] = 1$, then these two vertices have two links (or one bi-directional, if we allow it).

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Incidence matrix

- Each row is a vertex,
- Each column is an edge,
- Each value is a relation:
 - 0, if there is no relation between vertex and edge,
 - 1 if a vertex is a start of the edge,
 - 1 if a vertex is the end of the edge.



	e ₁	e ₂	e ₃	e ₄	e ₅	e ₆	e ₇
V ₁	-1	0	-1	0	0	0	1
V ₂	1	-1	0	0	0	-1	0
V ₃	0	0	0	0	1	1	-1
V ₄	0	1	1	-1	0	0	0
V ₅	0	0	0	1	-1	0	0

Properties of incidence matrix

- Much easier to add weights than in neighbourhood matrix.
- Each column must have one -1 and one 1 (integrity check).
- Number of 1s and -1s in rows → number of edges in and out.
- Finding neighbours is harder as we have to seek thru an entire row.
- All zeros in a row → insulated vertex.

	e ₁	e ₂	e ₃	e ₄	e ₅	e ₆	e ₇
V ₁	-1	0	-1	0	0	0	1
V ₂	1	-1	0	0	0	-1	0
V ₃	0	0	0	0	1	1	-1
V ₄	0	1	1	-1	0	0	0
V ₅	0	0	0	1	-1	0	0