# **Indexing Techniques for Vector Databases**

(practice & HW)

# 실습 목표 및 과제 소개

- IVF, HNSW 인덱스 동작 방식 이해
- Index 별 구성 요소 및 성능 평가 방식 이해

• 과제) IVF, HNSW + BQ 구현 (by python)

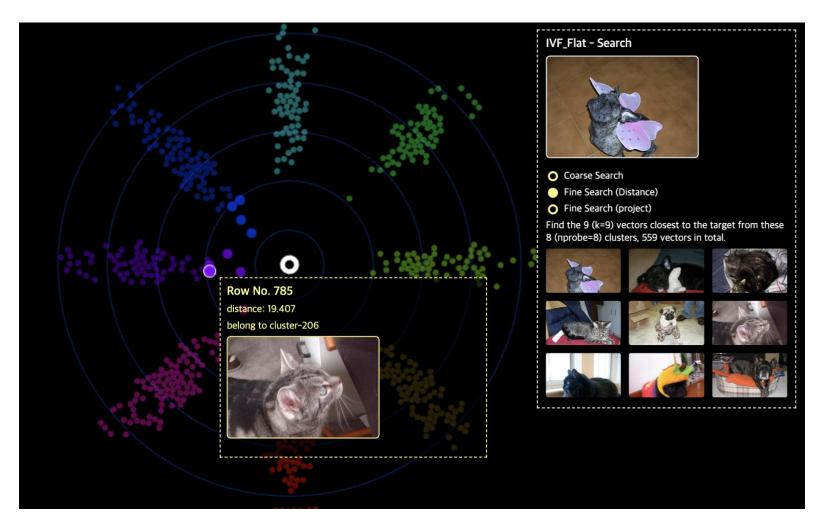
#### **IVF**

### IVF visualization (build)



• Clustering 진행 (region 수를 결정)

#### IVF visualization (search)



• 탐색할 probe 수 결정 (time <-> accuracy trade-off)

#### IVF python code

```
def __init__(self, distance_type, n_clusters):
    self.n_clusters = n_clusters
    self.distance_type = distance_type
    self.kmeans = KMeans(n_clusters=n_clusters, random_state=42)
    self.inverted_index = {i: [] for i in range(n_clusters)}
    self.data = None
```

- N\_cluster: cluster 개수
- Distance\_type: 사용할 distance measure
- Kmeans: clustering 시에 사용되는 모듈
- Inverted\_index: clustering 결과를 저장할 dictionary
- Data: 실제 벡터 데이터

#### IVF python code (build)

```
def fit(self, X):
    if self.distance_type == 'cosine':
        X = X / np.linalg.norm(X, axis=1, keepdims=True)

self.kmeans.fit(X)
labels = self.kmeans.labels_

for idx, label in enumerate(labels):
    self.inverted_index[label].append(idx)

self.data = X
```

• clustering 진행 후 inverted\_index에 저장

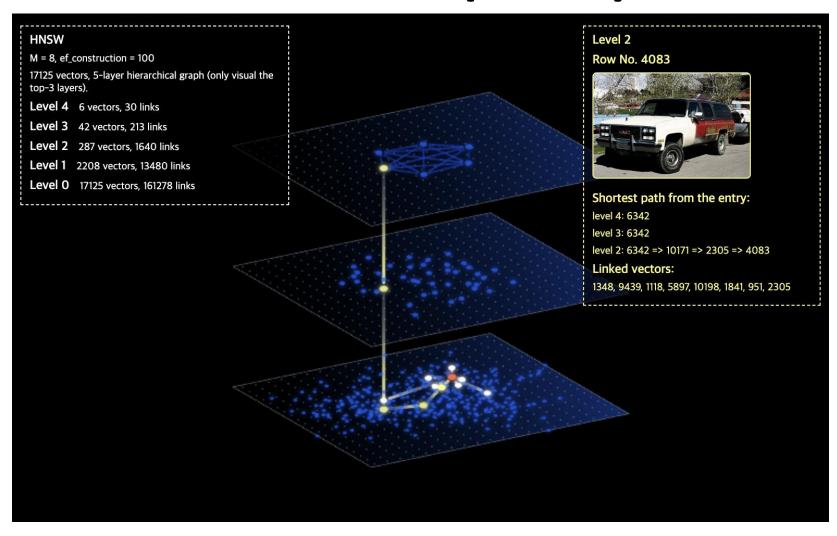
#### IVF python code (build)

```
def search(self, q, k, n_probes=1):
    assert n probes <= self.n clusters</pre>
    if self.distance type == "cosine":
        q = q / np.linalg.norm(q)
    cluster distances = self. compute distance(self.kmeans.cluster centers , q).flatten()
    probe_clusters = np.argpartition(cluster_distances, n_probes)[:n_probes]
    candidates = np.concatenate([self.inverted_index[cl] for cl in probe_clusters])
    if len(candidates) <= k:
        return candidates
    dists = self. compute distance(self.data[candidates], q).flatten()
    top_idx = np.argpartition(dists, k)[:k]
    top k = candidates[top idx]
    top dist = dists[top idx]
    return [(k, dist) for k, dist in zip(top_k, top_dist)]
```

- Cluster center와 query의 거리 계산
- n-probes 개의 근접 cluster 선택
- 해당 cluster 내에 있는 vector 중 상위 k개 반환

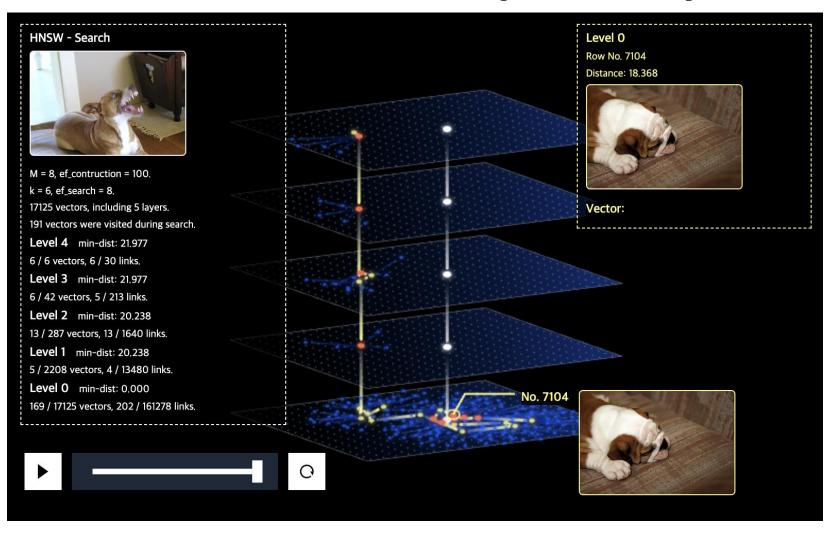
#### **HNSW**

#### **HNSW** visualization (build)



• 계층적 graph를 생성

#### **HNSW** visualization (search)



• Layer별로 search 진행

```
class HNSW(object):
    ### Algorithm 1: INSERT

def insert(self, q, efConstruction=None): "
    ### Algorithm 5: K-NN-SEARCH

def search(self, q, K=5, efSearch=20): "
    ### Algorithm 2: SEARCH-LAYER

def _search_layer(self, q, W, lc, ef): "
    ### Algorithm 3: SELECT-NEIGHBORS-SIMPLE

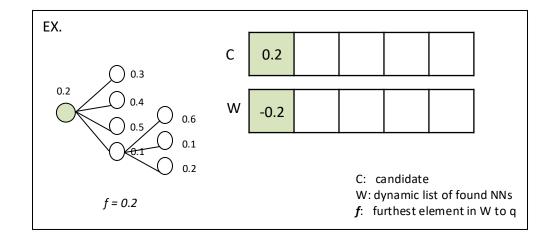
def _select(self, R, C, M, lc, heap=False): "
```

- 주요 메소드: insert, search
- 서브 알고리즘:
  - \_search\_layer: 주어진 계층에서 노드 탐색
  - \_select: 후보 노드 중 가장 가까운 이웃 연결

```
def __init__(self, distance_type, M=5, efConstruction=200, Mmax=None):
    if distance_type == "l2": "
        elif distance_type == "cosine": "
        else: "
        self.distance_func = distance_func
        self.vectorized_distance = self.vectorized_distance_
        self._M = M
        self._efConstruction = efConstruction
        self._level_mult = 1 / log2(M)
        self._level_mult = 1 / log2(M)
        self._enter_point = None
        self._data = []
```

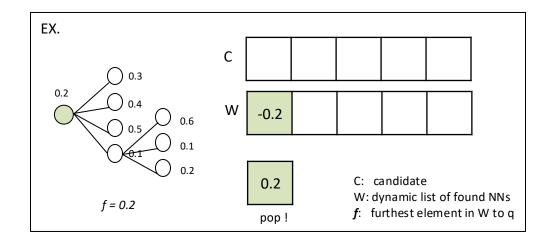
- M: 이웃 노드 수
- efConstruction: build 시 고려할 이웃 후보 수
  - Cf. efSearch: 검색 시 유지할 후보 노드 수
- level\_mult: 레벨 결정 시 사용되는 파라미터
- Graph, enter point, data -> 데이터 및 그래프 저장

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### Algorithm 2: SEARCH-LAYER
def search layer(self, q, W, lc, ef):
    vectorized distance = self.vectorized distance
   data = self.data
   # Step 1: Initialize candidate list and visited set
   C = [(-neg_dist, idx) for neg_dist, idx in W]
   heapify(C)
   heapify(W)
   visited = set(idx for _, idx in W)
    # Step 4-17: Explore neighbors until candidate list is exhausted
   while C:
        dist, c = heappop(C)
        furthest = -W[0][0]
        if dist > furthest:
            break
        neighbors = [e for e in lc[c] if e not in visited]
        visited.update(neighbors)
       dists = vectorized_distance(q, [data[e] for e in neighbors])
        for e, dist in zip(neighbors, dists):
            neg dist = -dist
            if len(W) < ef:
                heappush(C, (dist, e))
                heappush(W, (neg_dist, e))
                furthest = -W[0][0]
            elif dist < furthest:</pre>
                heappush(C, (dist, e))
                heapreplace(W, (neg dist, e))
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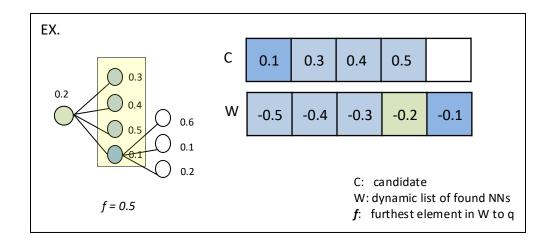
- 특정 layer에서 ef 개의 이웃 노드를 찾는 알고리즘
- C: 탐색 후보를 나타내는 min heap
- W: 검색 결과 후보를 나타내는 max heap
   최대 길이 ef
- Python의 경우 min heap만 제공, 따라서 W의 경우, distance를 음수값으로 저장하여 비교

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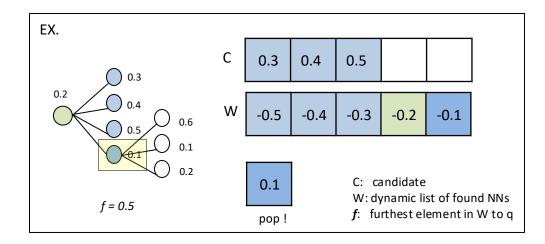
- C에서 pop (가장 가까운 이웃 후보 노드를 가져옴)
- 이를 w의 최댓값 (이미 뽑힌 노드 중 가장 먼 이웃)과 비교
- C에서 pop한게 더 먼 경우 종료
- 그렇지 않으면 이후 과정으로 넘어감

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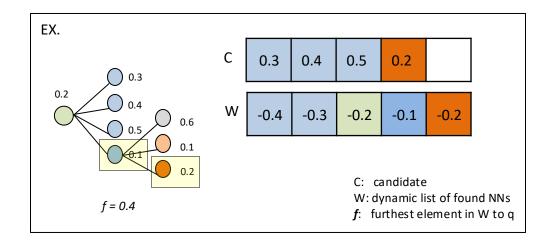


- Pop한 노드의 neighbor를 C, W에 추가
- W의 길이가 ef보다 작은 경우에 모든 노드를 추가

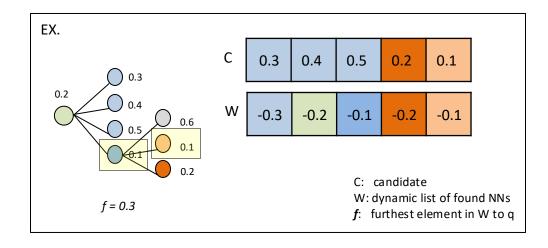
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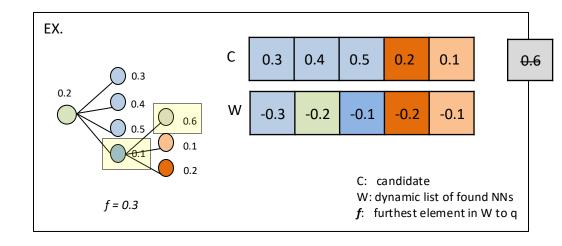
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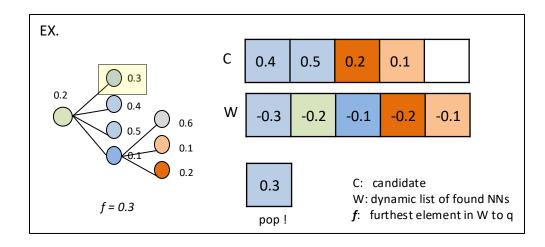
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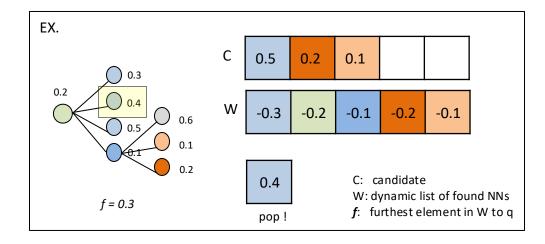
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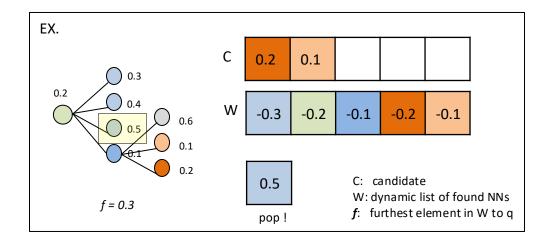
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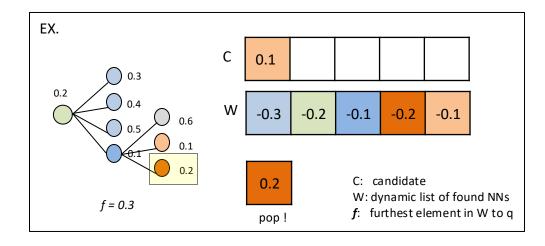
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```



```
### Algorithm 3: SELECT-NEIGHBORS-SIMPLE
def _select(self, R, C, M, lc, heap=False):

if not heap:
    idx, dist = C
    if len(R) < M:
        R[idx] = dist
    else:
        max_idx, max_dist = max(R.items(), key=itemgetter(1))
        if dist < max_dist:
            del R[max_idx]
            R[idx] = dist
    return

else:
    C = nlargest(M, C)
    R.update({idx: -neg_dist for neg_dist, idx in C})</pre>
```

- R: 이웃 노드 목록
- C: 추가할 노드 목록
- M: 최대 이웃 개수
- => R+C 중에서 M개를 선택하는 알고리즘

```
### Algorithm 1: INSERT
def insert(self, q, efConstruction=None):
   if efConstruction is None:
       efConstruction = self. efConstruction
   distance = self.distance func
   data = self.data
   graphs = self._graphs
    ep = self._enter_point
   M = self. M
   # line 4: determine level for the new element g
   l = int(-log2(random()) * self._level_mult) + 1
   idx = len(data)
   data.append(q)
   if ep is not None:
       neg_dist = -distance(q, data[ep])
       # distance(q, data[ep])
       # line 5-7: find the closest neighbor for levels above the insertion level
       for lc in reversed(graphs[l:]):
           neg_dist, ep = self._search_layer(q, [(neg_dist, ep)], lc, 1)[0]
       # line 8-17: insert q at the relevant levels; W is a candidate list
       laver0 = graphs[0]
       for lc in reversed(graphs[:l]):
           M layer = M if lc is not layer0 else self. Mmax
           # line 9: update W with the closest nodes found in the graph
            W = self._search_layer(q, [(neg_dist, ep)], lc, efConstruction)
            # line 10: insert the best neighbors for q at this layer
            lc[idx] = layer idx = {}
            self._select(layer_idx, W, M_layer, lc, heap=True)
            # line 11-13: insert bidirectional links to the new node
            for j, dist in layer_idx.items():
               self._select(lc[j], (idx, dist), M_layer, lc)
   # line 18: create empty graphs for all new levels
    for in range(len(graphs), l):
       graphs.append({idx: {}})
       self._enter_point = idx
```

• Level 결정

 Level보다 높은 layer에서는 제일 가까운 노드 하나만 찾아감

- Level 이하의 layer에서는 가까운 노드 efConstruction개를 찾음
- 그중 가장 가까운 M개를 골라, 본인과 이웃노드들의 이웃노드를 갱신

```
### Algorithm 5: K-NN-SEARCH
def search(self, q, K=5, efSearch=20):
    """Find the K points closest to g."""
    distance = self.distance_func
    graphs = self._graphs
    ep = self._enter_point
    if ep is None:
        raise ValueError("Empty graph")
    neg_dist = -distance(q, self.data[ep])
    # line 1-5: search from top layers down to the second level
    for lc in reversed(graphs[1:]):
        neg_dist, ep = self._search_layer(q, [(neg_dist, ep)], lc, 1) [0]
    # line 6: search with efSearch neighbors at the bottom level
    W = self._search_layer(q, [(neg_dist, ep)], graphs[0], efSearch)
    if K is not None:
        W = nlargest(K, W)
    else:
        W.sort(reverse=True)
    return [(idx, -md) for md, idx in W]
```

- Level 결정
- 맨 아래 layer를 제외, 가장 가까운 노드 하나만 찾아감
- 맨 마지막 layer에서 efSearch개를 탐색
- 탐색 결과에서 κ개를 선택

# 성능 평가 (practice.ipynb)

#### **Binary Quantization (practice.ipynb)**

# 과제 설명 ... (추가 예정)