S-Divergence-Based Internal Clustering Validation Index

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ABSTRACT

A clustering validation index (CVI) is employed to evaluate an algorithm's clustering results. Generally, CVI statistics can be split into three classes, namely internal, external, and relative cluster validations. Most of the existing internal CVIs were designed based on compactness (CM) and separation (SM). The distance between cluster centers is calculated by SM, whereas the CM measures the variance of the cluster. However, the SM between groups is not always captured accurately in highly overlapping classes. In this article, we devise a novel internal CVI that can be regarded as a complementary measure to the landscape of available internal CVIs. Initially, a database's clusters are modeled as a non-parametric density function estimated using kernel density estimation. Then the S-divergence (SD) and S-distance are introduced for measuring the SM and the CM, respectively. The SD is defined based on the concept of Hermitian positive definite matrices applied to density functions. The proposed internal CVI (PM) is the ratio of CM to SM. The PM outperforms the legacy measures presented in the literature on both superficial and realistic databases in various scenarios, according to empirical results from four popular clustering algorithms, including fuzzy k-means, spectral clustering, density peak clustering, and density-based spatial clustering applied to noisy data.

I. INTRODUCTION

CLUSTERING is an unsupervised methodology for analyzing a set of data objects by dividing them into subsets such that each group contains similar objects while dissimilar ones end up in different groups [1]–[5]. Thus, the objective of clustering is to mine the data to explore multi-dimensional obscure patterns and hidden structures in the data. Nowadays, clustering has received a great deal of attention among the community of researchers in the area of pattern recognition by the virtue of remarkable academic and commercial applications spanning over a wide range which includes identifying fake news [6], spam filtering [7], market segmentation [8], [9], classifying network traffic [10], detecting fraudulent or criminal activity [11], [12], cybersecurity [13], document analysis [14], drug discovery [15], information retrieval [16], and many more [17]–[22].

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A fundamental question in clustering is how to assess the "goodness" of the resulting clusters. The answer to this question is not obvious as it is difficult to devise criteria that determine the optimal partitioning of the data objects into clusters. Obtaining insights about the goodness of clusters using some visualization tools is not a feasible solution when the number of dimensions increases, as human eyes are not accustomed to higher-dimensional spaces. The process of assessing the performance of the clustering algorithm is referred to as cluster validation. According to the clustering validation procedure, the outcome of the clustering phase is validated quantitatively by a Clustering Validation Index (CVI). A CVI can be considered a function that, for a given clustering scheme and database, produces some value that represents the quality of the clustering scheme [23], [24]. In other words, a CVI provides some insight into the quality of grouping. Internal, external, and relative are the three main categories of CVIs. Internal CVIs rely only on the internal information of a given database. Unlike internal CVIs, external CVIs assess the "goodness" of a clustering structure based on provided class labels as external inputs [25]-[28]. On the other hand, relative CVIs evaluate the clustering

Keywords

Cluster Validity Index, Generalized Mean, K-nearest Neighbors, S-distance, S-divergence, Spectral Clustering, Symmetry Favored.

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(a) An instance of well separated clusters



Fig. 1. Distribution of three clusters.

results by changing the number of clusters. We mainly concentrate on internal CVIs in this work. The most intuitive notions for defining "good clustering" are cohesion/compactness (CM) and separation (SM). In simple words, when data objects in a cluster are in the vicinity of each other, the cluster is called a compact cluster. On the other hand, when neighboring clusters are possibly quite far from each other, then these clusters are easily identifiable and well separated. In other words, SM measures the distance between the centers of two clusters, whereas CM measures the variance within a cluster. Generally, geometric distance is used to compute SM. However, geometric distance can not always represent the SM efficiently, especially when two clusters are highly overlapping. Let us consider an example where three clusters, namely C_m , C_n , and C_p , are well separated (see Fig. 1(a)). As clusters are well separated, geometric distance can efficiently capture the dissimilarity between clusters. We may assume another scenario (see Fig. 1(b)), where clusters C_m , C_p , and C_p , are overlapping. In this case, the geometric distance between the centers of C_m and C_n is the same as the geometric distance between the centers of C_m and C_n . Thus, the dissimilarity between clusters can not be captured accurately using geometric distance. In [29], Cui et al. assumed that the data of a cluster were obtained from multivariate Gaussian distributions, and Jeffrey divergence (JD) was considered, as a distance measure for computing SM between clusters. The JD is not a valid distance measure because it does not abide by the metric property of triangle inequality [30]. In addition, the JD is not appropriate, while clusters are almost identical. Alternatively stated, a small change in clusters cannot be captured by JD. It encourages us to delve further in this direction by proposing an internal CVI based on the notion of S-divergence (SD), which can catch tiny variations in clusters since the cone is formed by the Hermitian positive definite matrices (HPDM). In addition, it fulfills all the properties of the distance metric [31]. Four well-known clustering techniques are employed to evaluate the performance of the proposed CVI on ten real-world and artificial databases. However, each cluster is modeled as a random variable using a non-parametric probability density function named kernel density estimation (KDE)

before the use of the proposed internal CVI. Among ten databases, some are well separated, a few are slightly overlapping, and the rest are highly overlapping. Moreover, noise is added in some databases to validate the efficacy of the proposed CVI. A comparative analysis is also performed to show the competitiveness of the proposed internal CVI in comparison with other CVIs.

The remaining article is structured as follows. After the introduction, in Section II, we examine several well-known internal CVIs. The proposed internal CVI is discussed further in Section III. Section IV provides the results of the experiments. At last, Section V concludes the work.

II. RELATED WORKS

A summary of some of the most popular internal CVIs is presented in this section. The CM reflects the average closeness or similarity of data points in all clusters. A value approaching 0 indicates good clustering [32]. The SM portrays the degree of separation between clusters [32]. A higher value of SM signifies better clustering. It is worth mentioning that other indexes, for example, root mean square standard deviation index (RMSSTDI) [33], root squared index (RSI) [33], and modified Hubert validity index (MHI) [34] perform on a different principle. Indeed, the RMSSTDI quantifies the homogeneity of the resultant clusters by calculating the square root of the aggregated variance of all the data objects. RSI determines the magnitude of difference between clusters using the ratio of the addition of the squares between-clusters to the total summation of the squares in the database. RMSSTDI, RSI, and MHI evaluate the difference betweenclusters by calculating the disagreements of groups of data objects in two parts. Furthermore, these indexes do not consider both CM and SM to validate the formed clusters. The Calinski-Harabasz index (CHI) computes the ratio of the sum of the average of between-clusters and of intra-cluster dispersion for all clusters [35]. A greater value of CHI demonstrates better partitions. CHI is usually fast to compute. Moreover, it is suitable for convex and well-separated clusters. On the other hand, it produces a low value for non-convex clusters. The Dunn validity index (DVI) calculates the SM of clusters over the CM of clusters [36]. Thus, a larger value of DVI suggests well-separated and compact clusters. However, the complexity of the DVI increases with the increase in the number of clusters, k. The Davies-Bouldin index (DBI) computes cluster overlapping using the ratio of the sum of intracluster spread to between-cluster distance [37]. A value adjacent to 0 illustrates better partitions. It computes the inherent attributes and quantities of a database. Moreover, it is limited to Euclidean space. The JD-based validity index (JI) is a ratio of CM to JD-based SM, [38]. JD determines the similarity between two probability distributions and is suitable for slightly-overlapping clusters. Thus, a value close to 0 is a sign of better partitions. However, JD falls short when the clusters are highly overlapping. The silhouette index (SI) measures how alike a data object is to its own cluster/cohesion/CM against other clusters/ SM [39]. A value near 1 signifies that the data object is well-suited to its cluster and does not match enough to neighboring clusters. A clustering configuration is appropriate when most data objects have a high value. SI is higher for well-separated and dense clusters. However, it is not suitable for non-convex clusters. Moreover, the computational complexity, $O(n^2 d \log(n))$, is high. I validity index (IVI) computes the CM and the SM using the maximum distance among data objects and centers of clusters [40]. Furthermore, the optimal number of clusters is calculated by maximizing the value of IVI. The Xie-Beni index (XBI) is defined using CM as the mean square distance among data objects and their cluster centers and the SM as the minimum square distance between the centers of clusters [41]. Optimal clusters exhibit a minimum value of XBI. The value of XBI reduces monotonically as the value of k increases. Furthermore, Bouguessa et al. [42] and Arbelaitz

TABLE I. A REVIEW OF SOME OF THE POPULAR INTERNAL CVIS

S. No.	Internal CVI	Notation	Expression	Range	Optimal value	Complexity
1	Root mean square standard deviation index	RMSSTDI	$\left\{ \frac{\sum_{i=1}^{k} \left. \sum_{c_j \in C_i} \left\ \mathbf{c}_j - \mathbf{v}_i \right\ ^2}{d \sum_{l=1}^{k} \left(\mathcal{C}_l - 1 \right)} \right\}^{1/2}$	[0, +∞]	elbow	O(nd)
2	Root squared index	RSI	$\frac{\sum_{\mathbf{c}\in DB} \ \mathbf{c} - \mathbf{v}\ ^2 - \sum_{i=1}^k \sum_{c_j \in C_i} \ \mathbf{c}_j - \mathbf{v}_i\ ^2}{\sum_{\mathbf{c}\in DB} \ \mathbf{c} - \mathbf{v}\ ^2}$	[0, 1]	elbow	O(nd)
3	Modified Hubert validity index	MHI	$\frac{2}{n(n-1)} \sum_{\mathbf{c}_j \in \mathcal{C}_i \text{ and } \mathbf{v}_i \in \mathcal{C}_i} \sum_{\mathbf{c}_q \in \mathcal{C}_r \text{ and } v_r \in \mathcal{C}_r} \text{ dist } (\mathbf{c}_j, \mathbf{c}_q) \text{ dist } (\mathbf{v}_i, \mathbf{v}_r)$	[0, +∞]	elbow	$O(n^2 d)$
4	Compactness measure	СМ	$\frac{1}{k}\sum_{i=1}^{k}\frac{1}{ C_i }\sum_{\mathbf{c}_j\in C_l}dist(\mathbf{c}_j,\mathbf{v}_l)$	[0, +∞]	Min	O(nd)
5	Separation measure	SM	$\frac{2}{k^2-k}\sum_{i=1}^k\sum_{p=i+1}^k\operatorname{dist}\left(\mathbf{v}_i,\mathbf{v}_p\right)$	[0, +∞]	Max	$O(k^2 d)$
6	Calinski-Harabasz index	CHI	$\sum_{i=1}^{k} \frac{ C_i \times \frac{dist(\mathbf{v}_i, \mathbf{v})}{(k-1)}}{\sum_{c_j \in C_i} \frac{dist(\mathbf{c}_j, \mathbf{v}_i)}{(n-k)}}$	[0, +∞]	Max	O(nd)
7	Dunn validity index	DVI	$\frac{\min_{1 \leq i \neq r \leq k} \left(\min_{\forall c_{c_0 \in C_0 \times c_q \in C_r}} \{ \text{dist} \left(\mathbf{c}_j, \mathbf{c}_q \right) \} \right)}{\forall c_{c_j} \in \mathcal{C}_{i,j}, \forall c_q \in C_r}$	[0, +∞]	Max	$O(n^2 d \log(n))$
8	Davies-Bouldin index	DBI	$\frac{1}{k} \sum_{i=1}^{k} \max_{\substack{r \neq i}} \left(\frac{\left \frac{1}{ c_i } \sum_{c_j \in c_i} \operatorname{dist} \left(\mathbf{c}_j, \mathbf{v}_i \right) + \frac{1}{ c_i } \sum_{c_j \in c_p} \operatorname{dist} \left(\mathbf{c}_j, \mathbf{v}_p \right)}{\operatorname{dist} \left(\mathbf{v}_i, \mathbf{v}_p \right)} \right)$	[0, +∞]	Min	$O(n^2 d \log(n))$
9	Jeffrey-divergence based validity index	JI	$\frac{\frac{1}{k}\sum_{i=1}^{k}\frac{1}{\Gamma_{i}}\sum_{c_{j}\in\mathcal{C}_{i}}dist(\mathbf{c}_{j},\mathbf{v}_{i})}{\frac{2}{k^{2}-k}\sum_{i=1}^{k}\sum_{p=i+1}^{k}JD(\mathbf{v}_{i},\mathbf{v}_{p})}$	[0, +∞]	Min	O(nd)
10	Silhouette index	SI	$\begin{split} &\frac{1}{n} \sum_{i=1}^{k} \sum_{\substack{c_j \in \mathcal{C}_i \\ max \{ sep (c_j, \mathbf{c}_q) - coh (\mathbf{c}_j, \mathbf{c}_l) \}}}{\max \{ sep (c_j, \mathbf{c}_q), coh (c_j, \mathbf{c}_l) \}}, \text{ where } coh (c_j, c_l) = \\ &\frac{1}{ \mathcal{C}_i } \sum_{\substack{c_j \in \mathcal{C}_i \\ c_j \in \mathcal{C}_i}} \text{ dist } (\mathbf{c}_j, \mathbf{c}_l) \text{ and } sep (c_j, \mathbf{c}_q) = \sum_{\substack{c_{rel} \text{ and } 1 < r < k}} \frac{1}{ \mathcal{C}_r } \sum_{\substack{c_q \in \mathcal{C}_r \\ c_q \in \mathcal{C}_r}} \text{ dist } (\mathbf{c}_j, \mathbf{c}_q) \end{split}$	[-1, 1]	Max	$O(n^2 d \log(n))$
11	I validity index	IVI	$\left(\frac{1}{k} \times \frac{\sum_{\boldsymbol{c} \in DB} \operatorname{dist} \left(\boldsymbol{c}, \boldsymbol{v} \right)}{\sum_{i=1}^{k} \sum_{\boldsymbol{c}, \boldsymbol{c} \in C_{i}} \operatorname{dist} \left(\boldsymbol{c}_{j}, \boldsymbol{v}_{i} \right)} \times \max_{1 \leq j, q \leq n} \left\{ \operatorname{dist} \left(\boldsymbol{c}_{j}, \boldsymbol{c}_{q} \right) \right\} \right)^{d}$	[0, +∞]	Max	$O(n^2 d \log(n))$
12	Xie-Beni index	XBI	$\sum_{i=1}^{k} \frac{\sum_{\mathbf{c}_{j} \in C_{i}} \operatorname{dist}^{2}\left(\mathbf{c}_{j}, \mathbf{v}_{i}\right)}{n \times \min_{\mathbf{c}_{j}, \mathbf{c}_{q} \neq c_{j}} \operatorname{dist}^{2}\left(\mathbf{c}_{j}, \mathbf{c}_{q}\right)}$	[0, +∞]	Min	$O(n^2 d \log(n))$

DB: Dataset, n: number of data objects in DB, v : center of DB, d: number of attributes, c: data objects of DB, k: number of clusters, C_i : i^{th} cluster, \mathbf{c}_j : j^{th} member of i^{th} cluster, \mathbf{v}_i : center of i^{th} cluster, v_i : center of C_i , dist (): distance function.

et al. [43] also worked to introduce indices based on Dunn variations and cohesion, which act well with noisy and overlapped clusters. Table I reports the definition, range, optimum value, and complexity of each of the above-discussed internal CVIs.

III. PROPOSED CVI

In this section, we examine and present some of the imperative properties of SD and propose a new internal CVI measure.

A. S-Divergence and Its Properties

Definition 1. SD presents a metric on the set of matrices A_r of size $\tau \times \tau$ [31]. The set A_r is a convex cone, on which SD is defined using Eq. 1.

$$D_{S}^{2}(A_{\tau}^{i}, A_{\tau}^{j}) = \log(\det(\frac{A_{\tau}^{i} + A_{\tau}^{j}}{2})) - \frac{1}{2}\log(\det(A_{\tau}^{i}A_{\tau}^{j}))$$
(1)

where det(.) denotes the determinant operation. D_s is a metric on the positive definite matrices (PDM) A_τ . Let ϕ_τ be a one-to-one function from $\Re^{\tau}_{+} \rightarrow A_{\tau}$. Now examine a vector $\mathbf{t} = \{t_1, t_2, ..., t_{\tau}\} \in \Re^{\tau}_{+}$ to generate PDM from a vector \mathbf{t} . SD is a divergence function on the cone of HPDM. A convex cone structure on the set of HPDM enables "geometric optimization", which enables us to resolve certain problems that may be non-convex in Euclidean space but convex in manifold space, or, offers efficient optimization. Thus, the divergence function on the cone of hpd matrices has empirical and computational advantages in many applications [44]. At this juncture, we shall demonstrate that the SD meets all the necessary characteristics for becoming a distance metric, which are given below:

Proposition 1. *Non-negativity*: $D_S^2(\phi_{\tau}(\mathbf{t}), \phi_{\tau}(\mathbf{u})) \ge 0$

Proof. The modified version of Eq. 1 is given below:

$$D_{S}^{2}(\phi_{\tau}(\mathbf{t}),\phi_{\tau}(\mathbf{u})) = \log(\det(\frac{\phi_{\tau}(\mathbf{t}) + \phi_{\tau}(\mathbf{u})}{2})) + \log(\frac{1}{\sqrt{\det(\phi_{\tau}(\mathbf{t})\phi_{\tau}(\mathbf{u}))}})$$
(2)

$$\Rightarrow D_{S}^{2}(\phi_{\tau}(\mathbf{t}),\phi_{\tau}(\mathbf{u})) = \log(\frac{\det(\frac{\phi_{\tau}(\mathbf{t})+\phi_{\tau}(\mathbf{u})}{2})}{\sqrt{\det(\phi_{\tau}(\mathbf{t})\phi_{\tau}(\mathbf{u}))}})$$
(3)

where $\frac{\det(\frac{\phi_{\tau}(t)+\phi_{\tau}(u)}{L})}{\sqrt{\det(\phi_{\tau}(t)\phi_{\tau}(u))}} \ge 0$ because determinant of the PDM is always positive and numerator will be greater than equal to denominator. $\therefore D_{S}^{2}(\phi_{\tau}(\mathbf{t}), \phi_{\tau}(\mathbf{u})) \ge 0$

Proposition 2. Equality:
$$D_S^2(\phi_{\tau}(\mathbf{t}), \phi_{\tau}(\mathbf{u})) = 0$$
 iff $\mathbf{t} = \mathbf{u}$

Proof. From proposition 1, we can write

$$D_{\mathcal{S}}^{2}(\phi_{\tau}(\mathbf{t}),\phi_{\tau}(\mathbf{u})) = \log(\frac{\det(\frac{\phi_{\tau}(\mathbf{t}) + \phi_{\tau}(\mathbf{u})}{2})}{\sqrt{\det(\phi_{\tau}(\mathbf{t})\phi_{\tau}(\mathbf{u}))}})$$

Now, if **t** and **u** are equal then **u** can be replaced by **t** in the above expression and the modified expression is

$$\begin{aligned} D_{S}^{2}(\phi_{\tau}(\mathbf{t}),\phi_{\tau}(\mathbf{t})) &= \log(\frac{\det(\frac{\varphi_{\tau}(\mathbf{t})+\varphi_{\tau}(\mathbf{t})}{2})}{\sqrt{\det(\phi_{\tau}(\mathbf{t})\phi_{\tau}(\mathbf{t}))}}) \Longrightarrow \log(1) = 0\\ \therefore D_{S}^{2}(\phi_{\tau}(\mathbf{t}),\phi_{\tau}(\mathbf{u})) &= 0 \text{ iff } \mathbf{t} = \mathbf{u}. \end{aligned}$$

Please note that we used the property that the determinant of the power of a matrix is equal to the determinant raised to that power, meaning in our case:

 $\det(\phi_{\tau}(\mathbf{t})\phi_{\tau}(\mathbf{t})) = \det(\phi_{\tau}(\mathbf{t}))^2$

Proposition 3. Symmetry: $D_S^2(\phi_\tau(\mathbf{t}),\phi_\tau(\mathbf{u})) = D_S^2(\phi_\tau(\mathbf{u}),\phi_\tau(\mathbf{t}))$

Proof. The SD amid **t** and **u** is denoted as follows:

 $D_{S}^{2}(\phi_{\tau}(\mathbf{t}),\phi_{\tau}(\mathbf{u})) = \log(\frac{\det(\frac{\phi_{\tau}(\mathbf{t})+\phi_{\tau}(\mathbf{u})}{2})}{\sqrt{\det(\phi_{\tau}(\mathbf{t})\phi_{\tau}(\mathbf{u}))}}) \text{ [as already noted in proposition } 1] = D_{S}^{2}(\phi_{\tau}(\mathbf{u}),\phi_{\tau}(\mathbf{t}))$ $\therefore D_S^2(\phi_\tau(\mathbf{t}), \phi_\tau(\mathbf{u})) = D_S^2(\phi_\tau(\mathbf{u}), \phi_\tau(\mathbf{t}))$

It implies SD also abides the symmetric metric property.

Proposition 4. Triangle Inequality: Suppose t, u, and z be three vectors. Then this proposition states, the sum of the lengths of any two sides viz., $D_{s}(\phi_{1}(\mathbf{t}), \phi_{1}(\mathbf{u}))$ and $D_{s}(\phi_{1}(\mathbf{u}), \phi_{1}(\mathbf{z}))$ of a triangle is greater than or equal to the length of the third side $D_s(\phi_{\tau}(t), \phi_{\tau}(z))$. Arithmetically, $D_{s}(\phi_{\tau}(\boldsymbol{t}),\phi_{\tau}(\boldsymbol{z})) \leq D_{s}(\phi_{\tau}(\boldsymbol{t}),\phi_{\tau}(\boldsymbol{u})) + D_{s}(\phi_{\tau}(\boldsymbol{u}),\phi_{\tau}(\boldsymbol{z})).$

Proof. Let **t**, **u**, and **z** be three vectors. Then $\phi_{z}(t)$, $\phi_{z}(u)$, $\phi_{z}(z) > 0$ and diagonal matrices.

Thus
$$D_S^2(\phi_{\tau}(\mathbf{t}), \phi_{\tau}(\mathbf{u})) = \sum_i D_S^2(t_i, u_i),$$

 $D_S^2(\phi_{\tau}(\mathbf{t}), \phi_{\tau}(\mathbf{z})) = \sum_i D_S^2(t_i, z_i), \text{ and}$
 $D_S^2(\phi_{\tau}(\mathbf{u}), \phi_{\tau}(\mathbf{z})) = \sum_i D_S^2(u_i, z_i)$
 $\therefore D_S^2(\phi_{\tau}(\mathbf{t}), \phi_{\tau}(\mathbf{z})) \le D_S^2(\phi_{\tau}(\mathbf{t}), \phi_{\tau}(\mathbf{u})) + D_S^2(\phi_{\tau}(\mathbf{u}), \phi_{\tau}(\mathbf{z}))$

Hence, it is showed that the SD is a metric.

B. Cluster Density Estimation

In this study, each cluster is modeled using a random variable characterized by a probability distribution. In practice, the underlying probability distribution of a random variable is not known in advance. Alternatively, the probability distribution of a random variable is estimated from the data objects or samples of a cluster. Therefore, each random variable is associated with a set of samples. We assume that samples are finite, independent, and identically distributed. Here, we adopt the well-known non-parametric probability estimation technique KDE to estimate the underlying distribution of the observations.

Let *M* be a random variable characterizing cluster C_m , where each sample, \mathbf{x} , is of *d*-dimensions. Then, the kernel function is obtained by multiplying the d number of Gaussian functions with bandwidth, h_l^M , where $1 \le l \le d$ and $d \ge 2$. Equation 4 is applied to estimate *M* [1], [2].

$$M(\mathbf{x}) = \frac{1}{|C_m|(2\pi)^{\frac{d}{2}} \prod_{l=1}^d h_l^m} \sum_{\mathbf{c}_j \in C_m} \prod_{l=1}^d e^{-\frac{(\mathbf{x}^l - \mathbf{c}_{j,l})^2}{2h_l^m}}$$
(4)

where $x \in D$, every cluster is defined in the same domain D and we also assume that the \mathcal{D} is a bounded range of values and c_i is a j^{th} member of i^{th} cluster or $c_i \in C_i$, h_l^M is the bandwidth of the $l^{th'}$ feature and it controls the smoothing of the Gaussian kernel function. The Sliverman approximation rule (Eq. 5) is considered to estimate h_1^M .

$$h_l^M = 1.06 \times \sigma_l |C_m|^{-\frac{1}{5}}$$
⁽⁵⁾

where σ_l denotes the standard deviation of C_m for the l^{th} feature.

C. S-Divergence Between Two Clusters

The SD between two clusters is stated as follows:

Definition 2. Let C_m and C_n be two clusters. The *M* and *N* are the two probability mass functions (PMFs) of C_m and C_n respectively as defined in Eq. 4 with finite or countably infinite values in a discrete domain, \mathcal{D} . The SD between C_m and C_n is computed by Eq. 6.

$$D_{S}^{2}(M,N) = \log(\det(\frac{\phi_{[c_{m}]}(M) + \phi_{[c_{m}]}(N)}{2})) - \frac{1}{2}\log(\det(\phi_{[c_{m}]}(M)\phi_{[c_{m}]}(N)))$$
(6)

where we assume that *M* has C_m samples $M = \{x_1, x_2, ..., x_{|C_m|}\}$ and PMF of every uncertain object is converted into diagonal matrix using $\phi_{|C|}$ () function as follows:

$$\phi_{|\mathcal{C}_{m}|}(M) = \begin{bmatrix} M(x_{1}) & 0 & \dots & 0 \\ 0 & M(x_{2}) & \dots & 0 \\ \vdots & \vdots & \dots & \vdots \\ 0 & 0 & \dots & M(x_{|\mathcal{C}_{m}|}) \end{bmatrix}$$
and
$$\phi_{|\mathcal{C}_{m}|}(N) = \begin{bmatrix} N(x_{1}) & 0 & \dots & 0 \\ 0 & N(x_{2}) & \dots & 0 \\ \vdots & \vdots & \dots & \vdots \\ 0 & 0 & \dots & N(x_{|\mathcal{C}_{m}|}) \end{bmatrix}.$$

Sometimes, it is needed to smooth a PMF of a data object thus the probability values become non-negative in a domain since SD consists of a logarithmic function as shown in Eq. 6. Thus, Eq. 7 is employed for normalizing [1].

$$N'(x) = \frac{N(x) + \beta}{1 + \beta |\mathcal{D}|}$$
(7)

where β is a constant and the value of β lies between an interval [0, 1]. The $|\mathcal{D}|$ signifies the number of possible values in \mathcal{D} . Furthermore, the sum of integral of N'(x) over the entire \mathcal{D} is 1. Equation 8 is utilized to estimate error in smoothing.

$$|N'(x) - N(x)| = \left|\frac{1 - N(x)\beta}{\frac{1}{\beta} + |\mathcal{D}|}\right| \in [0, \frac{max\{1, |1 - \mathcal{D}|\}}{\frac{1}{\beta} + |\mathcal{D}|}]$$
(8)

The value of β is assigned to 0.001 in this work. The $\phi_{|C_{n-1}}$ function is used to convert probability distributions to HPDM. The HPDM are manifolds, which are similar to non-positive curvature [31]. The HPDM cone does not come with a natural similarity function for a data object, although, it has computational and empirical advantages. Now, Eq. 6 is further simplified as follows:

$$\begin{split} D_{S}^{2}(M,N) &= \log(\frac{1}{2^{[C_{m}]}} \times \det(\phi_{|C_{m}|}(M) + \phi_{|C_{m}|}(N))) + \log(\frac{1}{\sqrt{\det(\phi_{|C_{m}|}(M)\phi_{|C_{m}|}(N))}}) = \\ \log(\frac{1}{2^{[C_{m}]}} \times \frac{(M(x_{1}) + N(x_{1}))(M(x_{2}) + N(x_{2})) \dots (M(x_{|C_{m}|}) + N(x_{|C_{m}|}))}{\sqrt{M(x_{1})M(x_{2}) \dots M(x_{|C_{m}|})N(x_{1})N(x_{2}) \dots g(x_{|C_{m}|})}}) \\ D_{S}^{2}(M,N) &= \log(\frac{1}{2^{|C_{m}|}} \times \frac{(M(x_{1}) + N(x_{1}))}{\sqrt{(M(x_{1})N(x_{1})}} \times \frac{(M(x_{2}) + N(x_{2}))}{\sqrt{(M(x_{1})N(x_{2})}} \times \dots \times \frac{(M(x_{|C_{m}|}) + N(x_{|C_{m}|}))}{\sqrt{(M(x_{|C_{m}|})N(x_{|C_{m}|})}}) \\ D_{S}^{2}(M,N) &= \log(\frac{1}{2^{|C_{m}|}}) + \log((\sqrt{\frac{M(x_{1})}{N(x_{1})}} + \sqrt{\frac{N(x_{1})}{N(x_{1})}})(\sqrt{\frac{M(x_{2})}{N(x_{2})}} + \sqrt{\frac{N(x_{2})}{M(x_{2})}}) \dots (\sqrt{\frac{M(x_{|C_{m}|})}{N(x_{|C_{m}|})}} + \sqrt{\frac{N(x_{|C_{m}|})}{M(x_{|C_{m}|})}}) \\ D_{S}^{2}(M,N) &= \log(\frac{1}{2^{|C_{m}|}}) + \log(\sqrt{\frac{M(x_{1})}{N(x_{1})}} + \log(\sqrt{\frac{M(x_{2})}{N(x_{2})}} + \dots + \log(\sqrt{\frac{M(x_{|C_{m}|})}{g(x_{|C_{m}|})}}) + \log(1 + \frac{N(x_{1})}{M(x_{1})}) \\ &\quad + \log(1 + \frac{N(x_{2})}{M(x_{2})}) + \dots + \log(1 + \frac{N(x_{|C_{m}|})}{M(x_{|C_{m}|})}) \\ D_{S}^{2}(M,N) &= \log(\frac{1}{2^{|C_{m}|}}) + \sum_{x \in \mathcal{D}} \log(\sqrt{\frac{M(x_{1})}{N(x)}}(1 + \frac{N(x)}{M(x)})) \end{split}$$

Finally, the SD between *M* and *N* is expressed as follows:

$$D_{S}^{2}(M,N) = \log(\frac{1}{2^{|C_{m}|}}) + \sum_{x \in \mathcal{D}} \log(\frac{|C_{n}| \prod_{l=1}^{d} h_{l}^{N} \sum_{c_{j} \in C_{m}} \prod_{l=1}^{d} e^{\frac{(x^{l} - c_{j})^{2}}{2h_{l}^{M^{2}}}}}{|C_{m}| \prod_{l=1}^{d} h_{l}^{M} \sum_{c_{j} \in C_{m}} \prod_{l=1}^{d} h_{l}^{M} \sum_{c_{j} \in C_{m}} \prod_{l=1}^{d} e^{\frac{(x^{l} - c_{j})^{2}}{2h_{l}^{M^{2}}}})} + \frac{|C_{m}| \prod_{l=1}^{d} h_{l}^{M} \sum_{c_{j} \in C_{m}} \prod_{l=1}^{d} e^{\frac{(x^{l} - c_{j})^{2}}{2h_{l}^{M^{2}}}})}{|C_{m}| \prod_{l=1}^{d} h_{l}^{N} \sum_{c_{j} \in C_{m}} \prod_{l=1}^{d} e^{\frac{(x^{l} - c_{j})^{2}}{2h_{l}^{M^{2}}}})}$$

D. The Proposed Internal CVI

The proposed internal CVI (PM) is based on CM and SM. So, the values of CM and SM need to be computed before calculating PM. The CM indicates the closeness or similarity of data objects in a cluster. Moreover, it is an average CM of all k clusters. The CM of every cluster, C_p is calculated by Eq. 9. It is an average of aggregated squared S-distance (SD) of a cluster data object c_i to its center v_p .

$$CM = \frac{1}{k} \sum_{i=1}^{k} \frac{1}{|C_i|} \sum_{\mathbf{c}_j \in C_i} D_{SD} \left(\mathbf{c}_j, \mathbf{v}_i \right)$$
(9)

where D_{SD} is the SD, which can be defined mathematically using Eq. 10.

Definition 3. define D_{SD} : $\mathfrak{R}^d_+ \times \mathfrak{R}^d_+ \to \mathfrak{R}_+ \cup \{0\}$ as

$$D_{SD}(\mathbf{c}_{j}, \mathbf{v}_{i}) = \sum_{l=1}^{d} \left[\log((\mathbf{c}_{j,l} + \mathbf{v}_{i,l})/2) - (\log(\mathbf{c}_{j,l}) + \log(\mathbf{v}_{i,l}))/2 \right]$$
(10)

Equation 10 shows a point-to-point distance measure labeled as the SD that is motivated by the SD. It is defined in the open cone of PDM. Moreover, Eq. 10 shows that if two data objects with the same Euclidean distance are close to the origin, then data objects will have a larger SD compared to when they are far from the origin. This property can be applied to find the properties of clusters with varying sizes and densities. Furthermore, SD is neither an f-divergence nor a Bregman divergence and is invariant under the Hadamard product [45].

The CM ranges from 0 to ∞ , where a low value is appropriate for a clustering configuration. The SM determines the magnitude of separation between clusters. The SD-based SM is calculated in this study by Eq. 11.

$$SM = \frac{2}{k(k-2)} \sum_{i=1}^{k} \sum_{j=i+1}^{k} D_{S}(M_{i}, M_{j})$$
(11)

where M_i and M_j are the PMFs of clusters C_i and C_j respectively. The SM lies in the interval $[0, \infty)$, where a high value implies good clustering. The PM is a ratio of the CM to the proposed SM, and it is estimated using Eq. 12.

$$PM = \frac{CM}{SM} \tag{12}$$

Good clustering is characterized by a low CM and a high SM of clusters. Therefore, a smaller value of *PM* is suitable for a clustering configuration. Sometimes, it is required to normalize the SM, and thus its value becomes non-zero in a domain since the zero value of SM will make an undefined value of the proposed index, PM, as shown in Eq. 12. Hence, Eq. 11 is further normalized.

$$SM = \frac{2}{k(k-2)} \sum_{i=1}^{k} \sum_{j=i+1}^{k} D_S \left(M_i, M_j \right) + \frac{\delta}{k^2}$$
(13)

where δ is a constant and the value of $\delta \rightarrow 0$, further estimated error in normalization is $\frac{\delta}{k^2}$ which is less significant in the possible range of SM. Normalized SM will be used throughout the paper to avoid an undefined value.

E. Complexity Analysis

The complexity associated with CM and SM is O(nd) and $O(k^2dE)$ respectively, where *E* is the number of steps to estimate the SD between two clusters. The complexity of PM is represented by $O(nd + k^2dE)$ since $n \ge d$ and $n \ge k$ is considered in this study. Thus the complexity of the proposed CVI is linear.

IV. Experimental Results and Discussion

A laptop Intel(R) Core(TM) i7-2620M CPU@2.70GHz and 4-GB RAM running on Windows 10 having a 64-bits Python 3.6.5 compiler are considered for this study. All the work is carried out in Spyder 3.2.8's Python development environment.

A. Description of Databases

A total of 10 databases of two classes, namely synthetic and realworld are considered in this work to prove the effectiveness of the PM over some of the most popular existing internal CVIs. Synthetic databases: Three databases, namely Blobs, Varied Distributed data, and Anisotropically Distributed Data, are created in this study. The title of the databases, the total number of data objects in each database, the total number of features in each data object, and the number of clusters are noted in Table II. The Blobs database is produced by an isotropic Gaussian function with three classes having 1500 data objects or samples and two features. The varied distributed data is produced with varied variance in the data and has 1500 samples with 3 classes in 2D space, whereas Anisotropicly distributed database is generated by transforming the data, which is Anisotropically distributed or aligned on a specific axis. This database also has 1500 samples, three classes, and two features. UCI and Kaggle repository databases: Seven popular realistic databases, viz., Digits, Iris, Wine, Avila, Shuttle, Breast Cancer, and Letter Recognition, are adopted from the UCI repository [46], [47]. The short description of each of these UCI databases is also reported in Table II. All the databases are renamed as DB, where i varies from 1 to 10.

TABLE II. DATASETS CHARACTERISTICS

5. No.	Datasets	No of data objects	No of features	Clusters
1	Varied distributed data (DB1)	1500	2	3
2	Anisotropicly distributed data (DB2)	1500	2	3
3	Blobs (DB3)	1500	2	3
4	Breast cancer database (DB4)	569	30	2
5	Iris database (DB5)	150	4	3
6	Wine database (DB6)	178	13	3
7	Avila database (DB7)	10430	10	12
8	Digits database (DB8)	1797	64	10
9	Letter recognition database (DB9)	20000	16	26
10	Shuttle database (DB10)	43500	9	7

B. Results and Comparison

A couple of experiments are conducted to prove the effectiveness of PM over some of the existing internal CVIs in different scenarios, which are as follows:

1. The Impact of Monotonicity

The first experiment aims to study the monotonicity behavior of three internal CVIs, namely RMSSTDI, RSI, and MHI. Three synthetic databases, namely DB_1 , DB_2 , and DB_3 are considered, where clusters are well-separated. Fig. 2 (a), (c), and (e) plot the datasets DB_1 , DB_2 , and DB_3 along the *x* and *y* axes on a 2D plane, respectively. Here, fuzzy k-means (FKM) is applied to the three databases mentioned above, and the values of RMSSTDI, RSI, and MHI are computed, which are labeled as FKM-RMSSTDI, FKM-RSI, and FKM-MHI, respectively. Fig. 2 (b), (d), and (f) show the values of FKM-RMSSTDI, FKM-RSI, and FKM-MHI, respectively, that are obtained by varying the number of



Fig. 2. DB,, DB, and DB, are plotted on the plane, different classes are shown with different colors and result of internal CVIs on database in the right.

clusters, k, from 2 to 29 as inputs because the datasets discussed in Table II have an actual number of clusters in the range of 2 to 26. The other information on the results is not pertinent to this experiment. The vertical axis of curves or graphs in Fig. 2 is scaled for better visualization or analysis. When the value of *k* increases then value of numerator in $RMSSTDI = \frac{\sum_{i=1}^{k}\sum_{i \in j} |k| - v_i||^2}{d(n-k)}$ will decrease. The value of (n - k) is regarded as a constant because $k \ll n$. Therefore, RMSSTDI decreases with an increase in the *k*-value in Fig. 2 (b), (d), and (f). Further, RSI specifics a ratio of between clusters sum of squares to the total sum of squares. Hence, RSI increases as the value of *k* increases, as shown in Fig. 2 (b), (d), and (f). Similarly, MHI increases as the value of *k*

increases, according to Fig. 2 (b), (d), and (f), because with an increase in k more pairs of distances are calculated. Furthermore, RMSSTDI is only based on CM, and RSI and MHI rely only on SM. According to the property of monotonicity, the curves of RMSSTDI, RSI, and MHI will be either downward or upward. It is quoted that the value of kis optimal at the "elbow" point, where a shift in the curve appears. Thus, the empirical results in Fig. 2 prove that the RMSSRDI, RSI, and MHI monotonically decrease or increase as the number of clusters, k, increases in the range from 2 to 29. However, the determination of a shift in the curve is rather a tedious and subjective task, thus the monotonicity is not discussed in the further sections.



Fig. 3. An analysis of internal CVIs on well-separated databases.

2. The Impact of Well-Separated Clusters

The aim of the 2^{nd} experiment is to determine the optimal value of k for the databases, where well-separated clusters are present. The steps involved in estimating the optimal value of k for the best partitions using internal CVIs are as follows:

- Step 1: Initialize a clustering algorithm before applying it to a database.
- Step 2: A set of parameters of the algorithm is fixed in order to achieve clustering results.
- Step 3: Calculate the corresponding internal CVIs after clustering.
- Step 4: Select the optimal value of internal CVIs for best partitions.

Here, the values of six internal CVIs viz., SI, CHI, DBI, DVI, JI, and PM are computed after applying FKM and spectral clustering (SC) [48] on three databases, namely DB₁, DB₂, and DB₂ and results are reported in Fig. 3 (a), (b), and (c) respectively. The FKM-SI, FKM-CHI, FKM-DBI, FKM-DVI, FKM-JI, and FKM-PM specify the values of SI, CHI, DBI, DVI, JI, and PM after executing FKM while SC-SI, SC-CHI, SC-DBI, SC-DVI, SC-JI, and SC-PM are employed to represent the values of SI, CHI, DBI, DVI, JI, and PM after applying SC. Fig. 3 displays the values of FKM-SI, FKM-CHI, FKM-DBI, FKM-DVI, FKM-JI, FKM-PM, SC-SI, SC-CHI, SC-DBI, SC-DVI, SC-JI, and SC-PM that are obtained by varying the value of k in the range of 2 to 29. The optimal values of CVIs labeled by a hexagon marker in Fig. 3 specify either maximum or minimum values, which demonstrate the actual values of k in the databases. It is clear from Fig. 3 (a) that SC-PM, FKM-PM, FKM-JI, SC-JI, SC-DVI, SC-CHI, SC-SI, FKM-CHI, and FKM-SI determine the optimal value of k, which is the same as the exact number of clusters in DB, Moreover, the remaining CVIs produce values of k, which are closer to the actual number of clusters. It is also observed from Fig. 3 (b) that the FKM-PM and FKM-JI compute the optimal number of clusters, which are equal to the real number of clusters in DB₂. Furthermore, FKM-SI, FKM-DBI, SC-SI, SC-DBI, SC-DVI, SC-PM, FKM-JI, and SC-JI are also in proximity to the optimal clusters. On the other hand, the remaining CVIs are not near-optimal results. Fig. 3 (c) shows the results of DB_2 and that FKM-SI, FKM-CHI, FKM-DBI, FKM-PM, SC-DBI, SC-PM, FKM-JI, SC-JI, and SC-CHI achieve the optimal value for the clusters.

3. The Impact of Slightly Overlapped Clusters

The third experiment aims to decide the optimal value of k for the databases, namely DB_4 , DB_5 , and DB_6 , where slightly overlapping clusters are present. However, principal component analysis is adopted in exploratory data analysis by transforming the data to a new coordinate system in the case of high-dimensional data and then plotting the first two principal components [49], [50]. The first two principal components of datasets DB_4 , DB_5 , and DB_6 are mapped on a 2D plane, which are displayed in Fig. 4 (a), (c), and (e), respectively. Here, slightly overlapping clusters are denoted by different colors. Again, the values of six internal CVIs, viz. SI, CHI, DBI, DVI, JI, and PM, are computed after applying FKM and SC on the three databases mentioned above, and the outcomes are noted in Fig. 4 (b), (d), and (f), respectively. Here, we run the clustering algorithms for different values of k in the range of 2 to 29. We can find out the exact values of k by considering the optimum values of the curves of the FKM-PM and SC-PM in most cases. Moreover, PM always helps to decide the exact value of k because of the use of non-linear similarity measures.

4. The Impact of Highly Overlapped Clusters

The focus of the fourth experiment is to estimate the optimal value of k for the databases, namely DB7, DB8, DB9, and DB10, where clusters are highly significant. The first two principal components of datasets DB_{7} , DB_{8} , DB_{9} , and DB_{10} are mapped on a 2D plane, which are displayed in Fig. 5 (a), (c), (e), and (g), respectively. Here, different colors are employed to represent clusters. Again, the values of six internal CVIs, viz. SI, CHI, DBI, DVI, JI, and PM, are calculated after applying FKM and SC on the four databases stated above, and the results are displayed in Fig. 5 (b), (d), (f), and (h), respectively. Here, both the clustering algorithms execute for different values of k in the range of 2 to 29. Focusing on the results, PM determines the optimal kfor DB, and DB, But FKM-DBI, FKM-DVI, SC-SI, and SC-PM compute a value close to it. Furthermore, FKM-PM, SC-PM, and SC-DVI find the optimal k for DB_0 , and FKM-DBI and FKM-DVI are not far from them. Finally, for DB_{10} , SC-PM and SC-DVI find the optimal k and FKM-DBI, FKM-DVI, FKM-PM, SC-DBI, and SC-JI compute a close value.

5. The Impact of Noise

The purpose of the 5th experiment is to determine how robust the proposed internal CVI named PM is against noisy features. First, noisy facets are included in the three well-separated databases, namely DB_1 , DB_2 , and DB_3 . Here, a noisy feature is produced by considering uniform random distribution in the limit of the length and size similar to features of the original database. The number of features will be doubled in a database after adding noisy features. The impact of noisy features is then analyzed in this study. Databases are shown in Fig. 6 (a), (c), and (e). Again, the values of six internal CVIs, viz. SI, CHI, DBI, DVI, JI, and PM, are estimated after applying FKM and SC to the three noisy databases presented above, and the results are portrayed in Fig. 6 (b), (d), and (f), respectively. Here, both the clustering algorithms execute for different values of k in the range of 2 to 29. It is clear from Fig. 6 that DBI and DVI are affected by noise and face difficulty while determining the optimum value of k. Further, the curve of CHI is close to the optimal number of clusters in the case of a noisy DB2 database. On the other hand, the optimum values of SI, JI, and PM are closer to the exact values of k.

We can conclude from the five experiments conducted above that the proposed internal CVI named PM successfully ascertains the optimal number of clusters for most databases. On the other hand,



Fig. 4. In the left first two principal components of the DB_{s} , DB_{s} , and DB_{a} are plotted on the plane, to display the first and second corresponding vectors of the data matrix along the axes, different classes are shown with different colors and result of internal CVIs on database in the right.

JI, SI, CHI, DBI, and DVI face difficulty while estimating the exact number of clusters due to various degrees of overlapping between clusters and noise in the databases.

6. The Comparative Analysis

Finally, the PM is compared with five popular internal CVIs, namely SI, CHI, DBI, DVI, and JI, after applying four clustering algorithms, viz. FKM, SC, Density-based Spatial Clustering of Applications with Noise (DBSCAN), and Density Peak Clustering (DPC) [51] on the ten databases mentioned in Section IV A. Here, FKM and SC take the exact number of clusters as inputs, whereas DBSCAN and DPC compute

the number of clusters automatically. The values of six internal CVIs, including the PM, are reported in Table III. The mean (μ) and standard deviation (σ) obtained by the four clustering algorithms of each CVI are also noted in the last column of Table III. The μ and σ of the PM are highlighted by bold characters. A smaller value of σ in percentage specifies well-separated and compact clusters. In other words, a smaller value of σ demonstrates that the clustering configuration is appropriate. It is clear from Table III that the PM consistently outperforms five considered internal CVIs on ten databases in different scenarios presented in Table IV. Therefore, PM can be a great choice while evaluating clustering results.



Fig. 5. In the left first two principal components of the DB7, DB8, DB9, and DB10 are plotted on the plane, to display the first and second corresponding vectors of the data matrix along the axes, different classes are shown with different colors and result of internal CVIs on database in the right.

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	0177	TT73 -		DB40 AN	DBC	1 = 0/
Dataset	CVI	FKM	SC	DBSCAN	DPC	$\mu \pm \sigma \%$
DB_1	SI	0.6468	0.61745	0.57755	0.62765	0.61736 ± 8.79207
	CHI	5451.4914	3810.65643	3792.7863	4756.62157	4452.88893 ± 18.04829
	DBI	0.56956	0.54442	0.8377	0.61122	0.64073 ± 20.94112
	DVI	0.00785	0.01372	0.02252	0.03266	0.01919 ± 56.37716
	П	37.00884	39.56956	45.42931	34.47878	39.12162 ± 11.98999
	PM	29.56956	30.33441	34.71643	28.57965	30.80001 + 4.72936
DB	SI	0.63551	0.48386	0.40725	0 50914	0 50894 + 18 63688
	CHI	2882 88154	2202 55251	5465.06704	2202 72272	4114.03521 + 22.78248
	DDI	3883.88130	0.60640	3403.90704	5805.75875	0(5221 + 1)(5(517))
	DBI	0.49246	0.68642	0./1943	0.71491	0.00331 ± 10.00317
	DVI	0.00899	0.0069	0.00897	0.00376	$0.00/16 \pm 34.4/394$
	Л	44.00376	41.71511	57.67286	43.40933	46.70027 ± 15.80087
	PM	23.48942	30.68531	28.71825	22.72414	26.40428 ± 14.78514
	SI	0.4863	0.42646	0.33207	0.46153	0.42659 ± 15.85287
	CHI	2011.98126	1601.38907	1413.97578	1481.83841	1627.29613 ± 16.46313
	DBI	0.73157	0.78999	0.84494	0.82526	0.79794 ± 6.23412
DB_{3}	DVI	0.00825	0.01911	0.01358	0.00881	0.01244 ± 40.60672
	Л	47.79319	43.01848	49.44894	44.48351	46.18603 ± 6.39381
	PM	35.83244	39.79319	40.84494	37.81437	38.57124 ± 5.74615
DP	SI	0.69726	0 50825	0 509	0.67526	0.59744 ± 17.23188
	CHI	1300 20823	1089 92944	1245 56763	1251 53446	122180994 + 852372
	DBI	0 5044	0.62032	0.60006	0 55185	0.57366 ± 9.87326
	DVI	0.01721	0.02732	0.00900	0.00148	0.01462 ± 41.09165
		0.01/31	0.00726	0.01246	0.02148	0.01403 ± 41.98105
	JI	60.5044	68.01731	74.07588	63.92288	66.6 ± 8.76054
	PM	43.51121	49.63143	48.60891	41.56075	45.82808 <u>+</u> 7.46948
DB_{5}	SI	0.55282	0.55432	0.68674	0.68105	0.61873 ± 12.16705
	CHI	561.62776	558.05804	502.82156	513.92455	534.10798 ± 8.69226
	DBI	0.66197	0.64325	0.37927	0.39431	0.51970 ± 29.59093
	DVI	0.09881	0.12181	0.338	0.07651	0.15901 ± 76.31654
	JI	31.65626	32.11279	43.19802	32.38334	34.83760 ± 16.02200
	РМ	12.6709	14.65626	15.38275	13.40429	14.02855 ± 5.63466
DR	SI	0.56448	0.57114	0.56067	0.56203	0.56458 ± 6.23471
	CHI	552 85171	561 81566	670 62599	708 08668	623.34501 ± 12.48562
	DBI	0 53573	0 53424	0.55357	0 54434	0.54197 ± 1.64643
	БИ	0.00007	0.01626	0.0374	0.02200	0.02751 + 35.91714
	л П	48.016.36	0.01020 E8 E2E72	0.0374 E1.6427E	40.6252	5105776 ± 901017
)1 D14	46.01626	36.33373	51.04575	49.0555	31.93770 ± 0.93017
	PM	27.53573	30.53424	31.55357	31.49413	30.27942 ± 0.82341
	SI	0.1937	0.12995	0.1385	0.11951	0.14542 ± 22.77166
	CHI	5285.5617	4519.20875	4333.76871	4212.35646	4587.72391 ± 10.50701
DP	DBI	1.12112	1.29988	1.01121	0.8937	1.08148 ± 15.96829
	DVI	0.00182	0.00529	0.00197	0.00194	0.00276 ± 61.38810
	JI	24.12043	28.12995	35.10793	27.69293	28.76281 ± 15.97742
	PM	8.12138	8.28694	7.57481	6.22372	7.55171 ± 12.39664
	SI		0.1785	0.18289	0.17863	0.18066 ± 9.87675
	CHI	169.36261	161.20475	162.1034	171.6	166.07133 ± 3.12864
	DBI	1.9	1.88899	1.89937	1.84913	1.89023 ± 1.63817
DB ₈	DVI	0.21933	0.26126	0.17384	0.19023	0.21117 ± 18.15176
	Π	42 87789	39 26069	49 92082	41 99865	4351451 ± 1043365
	PM	15 92192	18 79859	18 90038	15 83872	1736490 ± 134074
		0.1462	0.150	0.14712	0.120	0.14630 ± 6.95094
DB_{9}	CLII	140	0.132	0.14/13	0.137	137625764 ± 625207
		142	490	140	/10/	1570.25704 ± 0.35237
	DRI	1.6855	1.63312	1.64295	1.35005	$1.3/91 \pm 9.73410$
	DVI	0.04536	0.04307	0.04136	0.04036	0.04254 ± 5.14657
	JI	82.65005	96.04536	99.02207	94.98688	93.17609 ± 7.75123
	PM	59.6855	66.62	60.63995	56.65005	60.89963 ± 3.57705
DB ₁₀	SI	0.97878	0.96967	0.97987	0.58508	0.87835 ± 22.26525
10	CHI	15723.30982	14946.92039	12879.555	16331.2233	14970.25213 ± 10.05018
	DBI	0.34179	0.25082	0.3709	0.44054	0.35101 ± 22.39261
	DVI	0.13045	0.24059	0.04701	0.09064	0.12717 ± 65.21501
	П	61.68367	58.31793	64.39526	63.44635	61.96080 ± 4.31864
	PM	40.33581	39.25111	43.36991	41.43915	41.09900 ± 4.27706



Fig. 6. In the left noisy-databases are plotted on the plane, different classes are shown with different colors and result of internal CVIs on noisy-database in the right.

TABLE IV. THE OVERALL REVIEW OF SOME INTERNAL CVIS

V. CONCLUSION

Index	Well- separated	Slightly- separated	Highly- overlapped	Noise
SI	G	G	Х	A
CHI	G	G	Х	A
DVI	G	А	Х	X
DBI	А	G	Х	A
JI	G	G	А	A
PM	G	G	G	G

Internal CVIs are employed frequently in clustering to measure the goodness of the clustering algorithms without taking any external inputs. Most of the existing internal CVIs depend on CM and the geometric distance-based SM when computing the distance between cluster centers. The previous studies showed that such CVIs are not capable of producing accurate results, especially when the clusters of a database are highly overlapping. As a remedy, we introduce a new internal CVI, PM, using a modified CM and an updated SM based on the notion of SD. Moreover, SD is defined on the cone of HPDM and is

shown to have experimental and computational advantages over the other approaches in many applications. On the other hand, SD is a point-to-point distance measure that is motivated by the definition of SD. It is defined in the open cone of PDM. Initially, clusters of a database are modeled using density functions by applying a nonparametric kernel density estimation method. The PM is defined as the ratio of the modified CM to the updated SM. A smaller value of the PM indicates that the clustering configuration is appropriate. Empirical results illustrate that the PM is proficient in determining the exact number of clusters and the best partition for several superficial and realistic databases, including the database with arbitrary cluster shapes. The proposed internal CVI faces difficulty in ascertaining the optimal number of clusters when noisy features are included in a few databases. In addition, the proposed internal CVI works efficiently for databases having only numerical attributes. The latter two aspects deserve further study. In future work, SD may be explored to develop an external CVI.

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Conflict of interest

The authors declare no conflict of interest.

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