Lecture Notes in Networks and Systems 613

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Third Congress on Intelligent Systems

Proceedings of CIS 2022, Volume 2



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Cite this paper

Lanjewar, M.G., Parate, R.K., Wakodikar, R., Parab, J.S. (2023). Detection of Starch in Turmeric Using Machine Learning Methods. In: Kumar, S., Sharma, H., Balachandran, K., Kim, J.H., Bansal, J.C. (eds) Third Congress on Intelligent Systems. CIS 2022. Lecture Notes in Networks and Systems, vol 613. Springer, Singapore.

https://doi.org/10.1007/978-981-19-9379-4_10

Download citation

.RIS业 .ENW业 .BIB业

DOI Published Publisher Name
https://doi.org/10.1007/9 19 May 2023 Springer, Singapore

78-981-19-9379-4_10

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Print ISBN Online ISBN 978-981-19-9378-7 978-981-19-9379-4

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Abstract

Detecting adulterants in turmeric is necessary because turmeric is a vital food constituent that adds color and flavor. In this work, the pure turmeric powders were mixed with starch to produce distinct concentrations (0, 10, 20, and 30%) (w/w). The reflectance spectra of samples were taken by visible–NIR spectroscopy. Spectroscopy in the wavelength range 400–1700 nm was used to record reflectance spectra of pure turmeric and starch–contaminated samples. The recorded spectra were preprocessed using a Savitzky–Golay filter and a second derivative with poly order of 2. The preprocessed spectra are then standardized, which are used to train and validate ML models. Three ML models were employed for classification: logistic regression (LR), K–nearest neighbor (KNN), and support vector machines (SVM). The LR and KNN obtained 100% accuracy, precision, recall, and F1–score, while SVM obtained 90% accuracy, 92% precision, 94% recall, and 91% F1–score. The performance of these models was tested with the stratified five fold method. The KNN model obtained the highest average accuracy of 92%, which is excellent compared to the other models.