

**Machine Learning for Soil and Crop Management**  
**Professor Somsubhra Chakraborty**  
**Agricultural and Food Engineering Department**  
**Indian Institute of Technology Kharagpur**  
**Lecture 22**

**Diffuse Reflectance Spectroscopy: Basics and Applications for Crop and Soil (Contd.)**

Welcome friends to the second lecture of week 5 of NPTEL an online certification course of machine learning for soil and crop management.

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The slide features a central image of the Indian Institute of Technology Kharagpur building, framed by a green and blue geometric border. To the right, the NPTEL logo is visible at the top, followed by the text 'NPTEL ONLINE CERTIFICATION COURSES'. Below this, the course title 'Machine Learning for Soil and Crop Management' is listed, along with the professor's name 'Prof. Somsubhra Chakraborty' and his affiliation 'Agricultural and Food Engineering Department, Indian Institute of Technology Kharagpur'. The slide also specifies 'Week 5: DIFFUSE REFLECTANCE SPECTROSCOPY: BASICS AND APPLICATIONS FOR CROP AND SOIL' and 'LECTURE 22'.

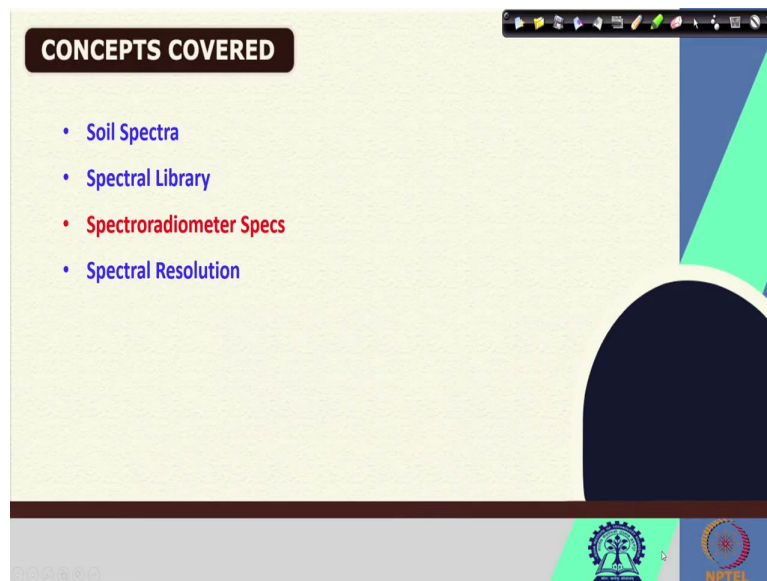
And in this week, our topic is Diffused Reflectance Spectroscopy: Basics and Applications for Crop and Soil. So, this is our second lecture and overall it is lecture number 22. So, in this lecture, we are going to discuss some of the important topics of diffuse reflectance spectroscopy and soil spectroscopy. In our previous lecture, we have covered some of the important aspects of what is panchromatic image, what is multispectral remote sensing, what is hyperspectral remote sensing, what is ultra-spectral spectra and then, what are the differences between their resolution and also we have focused on hyperspectral data cube and then we have started discussing the point spectroradiometers.

We have seen the difference between point spectroradiometers and imaging spectrometers, imaging spectrometers and imaging radiometers. We have seen some examples and then we have started discussing of the different aspects of soil spectroscopy we have defined what is called soil spectroscopy, we have seen the different zones like mid infrared spectra, then visible to near infrared spectroscopy, why mid infrared spectroscopy contains more information we have discussed because it contains the fundamental bands or fundamental

spectral features, which are arising from different soil minerals as well as from soil organic matter.

However, their overtones and combination bands are appearing in the visible to nearly infrared range. And that is why we required specialized spectral preprocessing as well as chemometric algorithms for resolving that complex pattern from the VisNIR-DRS spectrum. So, we have discussed all this.

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Now, in this lecture, we are going to start with the discussion of soil spectral. And then spectral library and then we are going to discuss some the important specification of the commercially available spectroradiometer and also we are going to discuss spectral resolution. Spectral resolution very important concept and why it is important we are going to discuss, what are the spectral resolution of spectroradiometers which we are also going to discuss.

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**KEYWORDS**

- Soil Spectra
- Spectral Library
- Spectral Resolution
- FWHM
- Sampling Interval

The slide features a presenter's video feed in the bottom right corner and logos for IIT Bombay and NPTEL at the bottom.

So, these are the some of the important keywords for this lecture, soil spectra, spectral liability, spectral resolution, FWHM or full width half maximum and then sampling interval. So, these terms will be discussed in this lecture.

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**HOW TO BUILD A SPECTRAL LIBRARY?**

The flowchart is divided into three main sections: BUILDING, USING, and a feedback loop.

- BUILDING:**
  - a) Legacy soil samples: Scan the soil → Outlier analysis → Outlier? → yes → Rescan → Scan the soil; no → Outlier? → yes → New targeted soil sampling → Collect soil samples → Scan the soil.
  - b) Define study area: Define a sampling design → Collect soil samples → Scan the soil → Outlier analysis → Outlier? → yes → Rescan → Scan the soil; no → Outlier? → yes → New targeted soil sampling → Collect soil samples → Scan the soil.
- USING:** Collect soil samples → Scan the soil → Does it fit in the spectral domain of the library? → no → Laboratory analysis → Predict soil properties; yes → Predict soil properties.
- Feedback Loop:** Predict soil properties → Measure laboratory accuracy → Laboratory analysis → Select subsample for laboratory analysis → Multivariate calibration → back to BUILDING.

The slide includes a presenter's video feed and logos for IIT Bombay and NPTEL at the bottom.

So, let us start with the building a spectral library. Now, it is when we talk about soil spectroscopy, it is very much important for to build a spectral library focusing on a study focusing on some soil properties focusing on some area region for building the future models as well as we can also utilize the legacy soil samples which have already collected before by different types of sampling procedures. So, that those are called the legacy soil samples.

So, which you have already measured in the lab and the properties have been already digitized and listed. So, these are known as the legacy soil samples. So, these legacy soil samples also can be used for developing of spectral library it is very much important for now, nowadays that you should build a spectral library for focusing on different types of soil covering a wide variation of soil properties. So, that you can develop the spectral algorithms for each individual region specific calibration model because there is there are certain problems while developing an universal model.

So, it has been recommended by several scientists that you should develop a region specific calibration model, spectral calibration model and for that you need to develop a spectral library. So, this spectral library can be also used for future matching of any soil spectra. Suppose, we have a large spectral library, we can capture, the when once you create this large spectral liability, which will be combining both the spectral information as well as the soil properties.

So, this spectral library can be in near future if we have developed a large spectral library combining both different soil properties as well as the spectral information then, you can use this library for a matching of a spectral matching for from an unknown soil samples which you can collect from any area of the world. So, the question comes how to build this spectral library? So, let us see the step by step which and this flowchart was given by the Viscarra Rossel and McBratney in the year 2008, which shows the step by step method of development of spectral library.

So, if we can start with the, there are 2 different part, one is building the spectral library, and then using the spectral library. So, when you concerned about the building the spectral libraries, we will start with the legacy soil samples, what are the legacy? So, we start with the legacy soil samples which have been already collected before. So, we scan the soil samples and then we go for the outlier analysis. Is there any outlier spectra or not? So, if there is an outlier, then if there is an outlier, then we should rescan the sample and then again check whether that is outlier or not.

If it is, if after the scanning if you see that there is no, that was a mistake and it was not actually an outlier then of course, we can go with the multivariate calibration using chemometric algorithm. However, if after rescanning, we still find that this is an outlier, then we should go with the new targeted soil sampling, because that may represent a new type of

soil and we should do the sampling of the same type of soil for building or augmenting the spectral library. Now, this is one of the way.

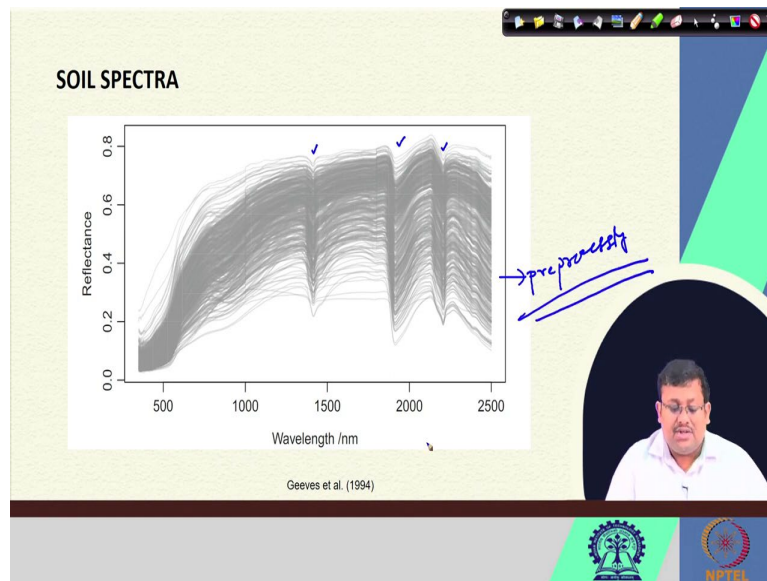
Also we can define a study area and then we can define a sampling design for collecting the samples from the study area. Once we define the study area define the sampling design, we can collect the soil samples. After collecting the soil samples, we can scan the soil samples, after scanning the soil samples, we can do the outlier analysis just like we have done in case of legacy soil data. So, if it is no if there is no outlier of course, we can select the sub sample for laboratory analysis and then we can develop the multivariate calibration as we have done in case of legacy soil samples.

However, if there is an outlier, then we should rescan the soil and then if there is of course, if the outlier was detected by mistake, then we can go back to this pathway. However, if it is still outlier, then we should go for the laboratory analysis and measure laboratory accuracy that will be an important issue. So, our laboratory analysis may be defective. So, we can check the laboratory analysis either way, or we can go with the new targeted soil sampling just we have done in case of legacy data.

So, this is about how to build a spectral library. However, how to use the specter library, once we develop this multivariate calibration model using chemometric algorithms that we can utilize for predicting the soil properties from unknown samples. So, we can collect the soil samples, we can scan the soil samples, and after scanning the soil samples, we will see does it fit in the spectral domain of the library or not by spectral matching. So, we can see whether that scan about the soil does fit in this spectral domain of the library or not.

So, if it is yes, then we predict the soil properties based on this multivariate calibration. If it is the answer is no, then we go for the laboratory analysis. So, this is how we generate and use the spectral library is very precisely given in this publication. So, those who are interested, I would request it, I would request you to go and see this publication for more important information.

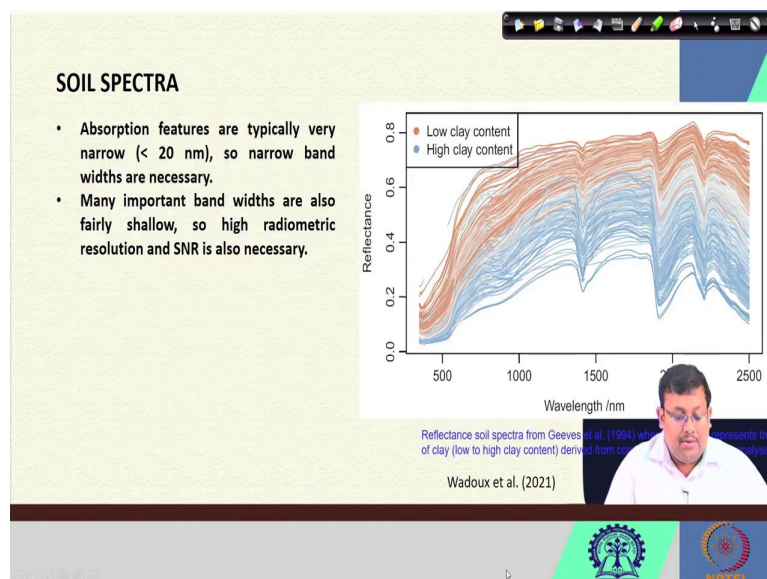
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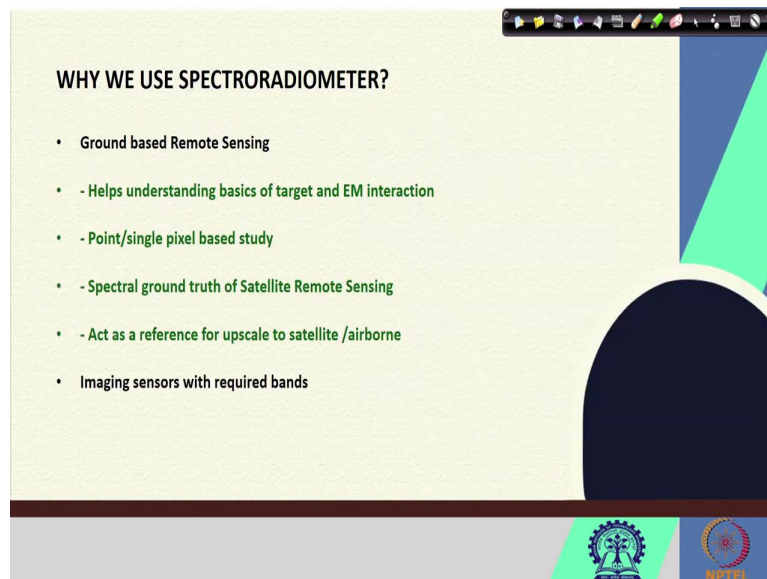


So, this plot shows the soil spectra, typical soil spectra of several soil samples, hundreds of soil samples and you can clearly see that the spectral features which are arising at 1450 nanometer and then 1900 around 1900 nanometer, 1910 nanometer and then the spectral features which are arising at around 2200 nanometer and so on so forth.

These are the broad spectra of the broad reflectance spectrum. However, this should be coupled with the spectral preprocessing to get more high signal to noise ratio. So, we will see that spectral preprocessing in our coming lectures.

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So, these are soil spectra, so, if we define the soil spectra so if you divide that soil spectra, which we have seen in our last slide into low clay content and high clay content, you can see the clear difference that low clay content soil sand clay, a high clay content soil in the spectral differences can be easily visible. Of course, because of high clay content, the concentration; high clay concentration, the soil will be much more darker and the reflection will be comparatively less than those which, for those soil which are having low clay content.

So, it is quite expected that the low clay content soil will have higher reflectance values than that of high clay content soil because of color difference. So, one thing you should understand that there is absorption features which are detected by the spectroradiometer should be are generally very narrow which are less than 20 nanometer. So, narrow band widths are necessary. So, we should select the spectroradiometer in such a fashion that each should capture these narrow spectral features.

So, that is why multispectral remote sensing or multispectral sensor cannot detect these minor spectral features which are less than 20 nanometer. So, that is why we require these hyperspectral point spectroradiometers which can which are having the narrow bandwidth and this narrow bandwidth is it is capable of capturing these capable of capturing these small spectral features. So, many important bandwidths are also fairly shallow. So, high radiometric resolution and SNR is also necessary.

So, not only the low bandwidth, but also high radiometric resolution and also signal to noise ratio is also necessary. Signal to noise ratio, high signal to noise ratio is important because we want to remove that noise and we want to keep the signal so, that or remove or reduce the

signal noise, so, that the signal to noise ratio can be maximized. So, that we can capture the targeted analyte more easily from the spectrum instead of a spectral curve which is having low SNR. So, low SNR is not preferable, high SNR is always preferable for prediction of a target analyte.

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**WHY WE USE SPECTRORADIOMETER?**

- Ground based Remote Sensing
- - Helps understanding basics of target and EM interaction
- - Point/single pixel based study
- - Spectral ground truth of Satellite Remote Sensing
- - Act as a reference for upscale to satellite /airborne
- Imaging sensors with required bands

So, the next question comes why we use spectroradiometer. So, generally, we use the spectroradiometer as far as the soil spectroscopy is concerned we generally use for ground based remote sensing for ground truthing whatever we can see from the satellite or airborne sensors, whether that reflects actually the surface features or not for that we need some ground truthing and for these ground truthing we use this point spectroradiometers. So, basically this spectrophotometer helps us for understanding the basics of the target and electromagnetic radiation interaction between this target and this radiation.

And also it helps in as I have told you it is a point spectrophotometer it helps in the point and singles pixel based study and then spectral ground truth of satellite remote sensing it is used for spectral ground truth of the satellite remote sensing we require and also we also require as a reference for upscale to satellite airborne. If we can develop any spectral algorithm using this point spectral radiometer, we can upscale it into the satellite and airborne sensor it is possible.

So, that in future by taking the hyperspectral image from the satellite platform or airborne platform, we can quantitatively measure the components within the image. So, this is why we use the spectroradiometer as far as the soil use is concerned. And in for in case of crop, we



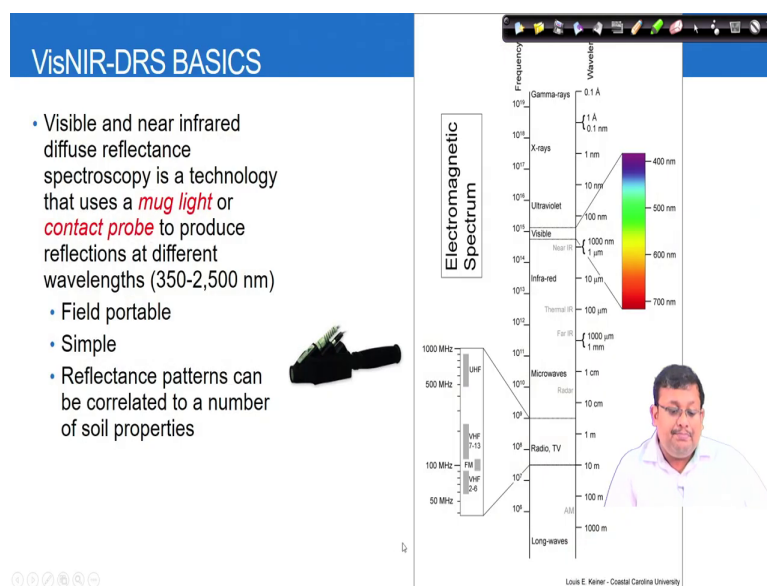
generally use a spectrophotometer for detecting the stress of the crop which I will discuss in our upcoming lectures also, so, for example, salinity stress and also we can generate some kind of spectral features or spectral index and we can calculate the spectral index also using the spectral features.

And also some time we use the spectral features or spectroradiometer for the, for capturing the environmental quality by seeing the phenotype, by seeing the spectral feature changes in the crop, which I am going to also discuss in our coming lectures. So, these are some of the basic requirements for using the spectrophotometer.

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### VisNIR-DRS BASICS

- Visible and near infrared diffuse reflectance spectroscopy is a technology that uses a *mug light* or *contact probe* to produce reflections at different wavelengths (350-2,500 nm)
- Field portable
- Simple
- Reflectance patterns can be correlated to a number of soil properties



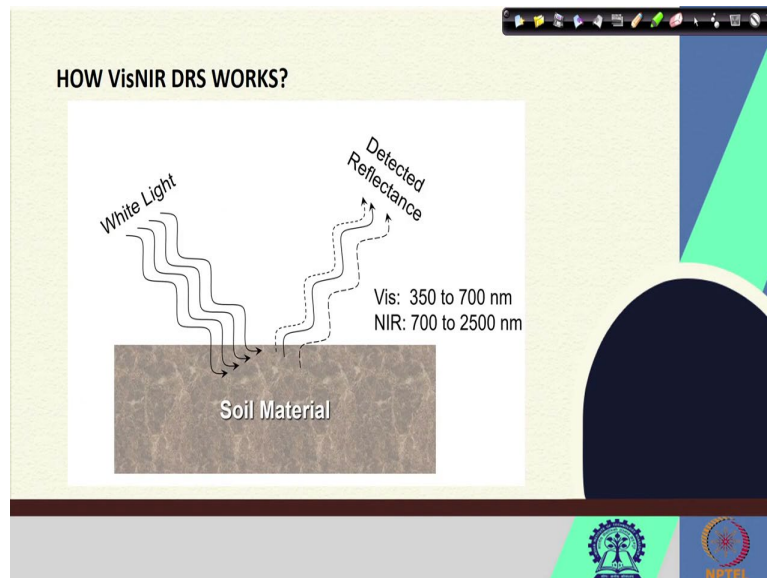
So, let us see the basics of VisNIR-DRS or visible near infrared diffuse reflectance spectroscopy. So, this visible and near infrared diffuse reflectance spectroscopy is a technology that uses a contact probe or mug light. So, this is called a contact probe, which is generally connected through the original spectroradiometer, it has its own light source and also it has a detector.

So, these produce when this light is coming from this window and reflects from any surface let us consider it is a soil, the soil will reflect some amount of radiation and this is reflected radiation will be captured from 350 nanometer to 2500 nanometer using the detector which is present inside this.

So, this is called the contact probe, there is are some similar type of fore-optics or similar types of fore-optics are there and attachments are there which we are also going to discuss. So, this is how these and once it captures this reflected energy, it transferred this reflected

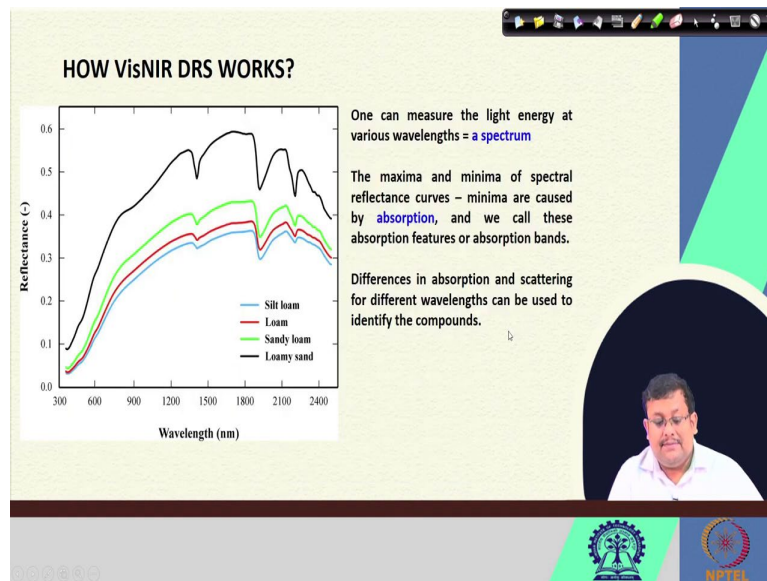
energy directly to the spectroradiometer through fiber optic cable. So, this is how this works and it is field portable, it is very simple to you handle and also the reflectance pattern can be correlated to a number of soil properties which we are going to discuss. So, this VisNIR-DRS basically captures the visible to near infrared range, which generally varies from the around 350 to 2500 nanometer region.

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So, we know that when there is a white light is coming and interacting with the soil material, some amount will be detected. So, these VisNIR-DRS there is basically captures this 350 to 2500 nanometer wavelength range in by their detector and measured the reflectance values or any other values like absorbance values using these detectors in the range of 350 to 2500 nanometer that means, it captures. Both 350 to 700 nanometers that is visible range and 700 nanometer to 2500 nanometers that is the near infrared region. So, visible to near infrared range, it can totally capture and then it can measure using this we know detected reflectance it predict soil properties.

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So, you know that one can measure the light energy at various wavelengths which is known as the spectrum and this is the typical soil spectrum with these are different 4 different types of soil silt loam, loam, sandy loam, loamy sand, and we can see that different types of spectral features are arising. So, the maxima and minima of the spectral reflectance curves, minima are caused by absorption. So, these are the minima. So, these are caused by the absorption.

And we call them the absorption features or absorption bands and the difference in absorption and scattering for different wavelengths can be used to identify the compounds. So, of course, the higher concentration of these compounds will result in higher absorption and lower concentration will result in lower absorption. So, we can, using the chemometric algorithm, if we capture this spectral variation through a quantitative way, we can predict the variation of certain soil parameters. So, this is how these VisNIR-DRS works.

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PSR+ 3500 Specifications		Performance Specifications		
Spectral Range	350-2500nm	Standard Res	Hi-Res	Hi-Res HD
Spectral Resolution-FWHM (Full Width at Half Maximum)	2.8nm @ 700nm 8nm @ 1500nm 6nm @ 2100nm	Resolution VNIR @ 700 nm	3 nm	3 nm
SI Detector	512 element Si photodiode array (350-1000nm)	Resolution SWIR @ 1400 & 2100 nm	8 nm	6 nm
InGaAs Detectors (cooled)	256 element extended wavelength photodiode array (970-1910nm) 256 element extended wavelength photodiode array (1900-2500nm)	Spectral Sampling (Bandwidth) VNIR	2.2 nm	2.2 nm
FOV Options (direct mount)	4°, 8°, or 14° lens, 25° fiber optic diffuser, integratinal sphere	Spectral Sampling (Bandwidth) SWIR	2.2 nm	2.2 nm
Fiber Mount Options	1, 2, 3, 4, 8 and 10° lenses	Scanning time	milliseconds	milliseconds
Noise Equivalence Radiance (4° lens)	0.5x10 <sup>-4</sup> W/cm <sup>2</sup> /nm/sr @400nm 0.8x10 <sup>-4</sup> W/cm <sup>2</sup> /nm/sr @1500nm 1.0x10 <sup>-4</sup> W/cm <sup>2</sup> /nm/sr @2100nm	NEΔ (Noise Equivalent Radiance)- VNIR @ 700 nm	1.0 x 10 <sup>-4</sup> W/cm <sup>2</sup> /nm/sr	1.0 x 10 <sup>-4</sup> W/cm <sup>2</sup> /nm/sr
Max Radiance @ 700nm (4° lens)	1.5x10 <sup>-4</sup> W/cm <sup>2</sup> /nm/sr	NEΔ- SWIR 1 @ 1400 nm	1.2 x 10 <sup>-4</sup> W/cm <sup>2</sup> /nm/sr	1.4 x 10 <sup>-4</sup> W/cm <sup>2</sup> /nm/sr
Minimum Scan Speed	100 milliseconds	NEΔ- SWIR 2 @ 2100 nm	1.9 x 10 <sup>-4</sup> W/cm <sup>2</sup> /nm/sr	2.2 x 10 <sup>-4</sup> W/cm <sup>2</sup> /nm/sr
Wavelength Reproducibility	0.1nm	Wavelength reproducibility	0.1 nm	0.1 nm
Wavelength Accuracy	±0.5 bandwidth	Wavelength accuracy	0.5 nm average error of wavelength calibration fit, Wavelength accuracy < 1 nm for any one line	0.5 nm
Communications Interface	USB or Class Bluetooth- laptop or PDA compatible	Maximum radiance- VNIR	2x Solar	2x Solar
Size	8.5" x 11.5" x 3.25"	Maximum radiance- SWIR	10x Solar	10x Solar
Tripod Mounting	2 each 1/4-20 mounting holes provided	Data collection speed	2 spectra per second	2 spectra per second
Weight	7.6 lbs (3.5 kg)	Channels	205	205
Batteries	Lithium ion, 7.4V, 7200mAh, 400mg/battery	SWIR (150-1000 nm) detector	512 element NIR enhanced silicon array	512 element NIR enhanced silicon array
Battery Operation	Rechargeable battery, typically up to 4 hour operation, battery (2 provided)	SWIR 1 (1000-1800 nm) detector	Graded InGaAs Photodiode, 2 Stage TE Cooled	Graded InGaAs Photodiode, 2 Stage TE Cooled
On Board Memory	Storage of 1000 spectra	SWIR 2 (1800-2500 nm) detector	Graded InGaAs Photodiode, 2 Stage TE Cooled	Graded InGaAs Photodiode, 2 Stage TE Cooled
		Input	1.5m fiber optic (25° field of view) optional fore optics available	1.5m fiber optic (25° field of view) optional fore optics available
		Weight	5.44 kg (12 lbs.)	5.44 kg (12 lbs.)
		Calibrations	Wavelength, absolute reflectance, radiance, irradiance. All calibrations are NIST traceable. Pushbutton calibrations are optional.	Wavelength, absolute reflectance, radiance, irradiance. All calibrations are NIST traceable. Pushbutton calibrations are optional.
		Instrument Controller	Dell Latitude 5480 or other Windows 10 compatible laptop	Dell Latitude 5480 or other Windows 10 compatible laptop
		Warranty	One year full warranty including expert customer support	One year full warranty including expert customer support
		Storage temperature (°C)	-15-45	-15-45
		Operational temperature range (°C)	0-40	0-40

So, these are the specification of the spectroradiometers you can see there are I have compared two most widely used spectroradiometers, one is PSR 3500 and then FieldSpec spectroradiometer. So, you can see both of them are having the spectral range of 350 to 2500 nanometer and the spectral resolution generally varies from one region to another region. So, you can see 2.8 nanometer at 700 nanometer, 8 nanometer 1500 nanometers, 6 nanometer at 2100 nanometer in case of PSR 3500. In case of FieldSpec, it is 3 nanometer 8 resolution at 1400; 2100 nanometer it is 8 nanometer.

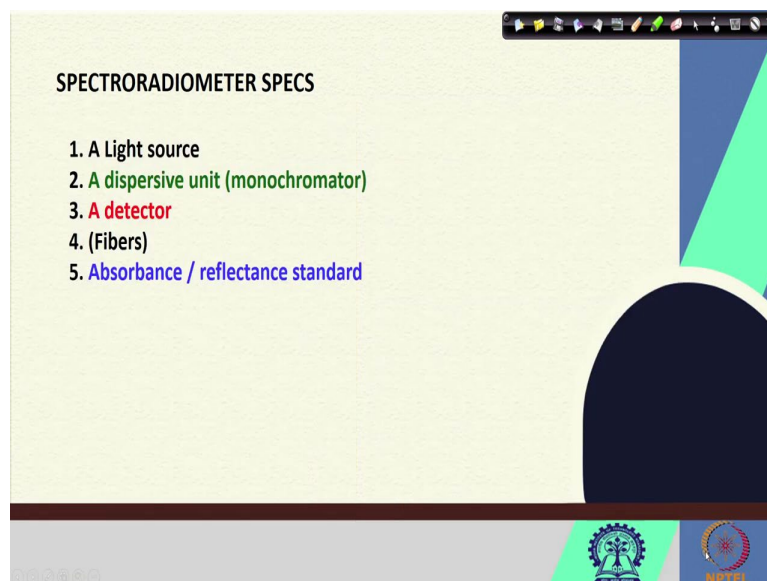
So, in case of the in the VNIR range, which is a, which is they called it is the given this specific name is given by the manufacturer which from the 350 to 1000 nanometer range, the detector is mainly 512 element silicon photodiode array which we are going to discuss. So, and also from the 970 to 1910 nanometer there are indium, gallium, arsenide detectors and also the 1900 nanometer to 2500 nanometer this is also indium, gallium, arsenide detector.

So, then there are other options like what are the field of views like 4 degree, 8 degree, 14 degree lens, 25 degree fiber optic diffuser integrating spear. So, these are different types of FOV options are there, what are the FOVs we are going to discuss also. So, here you can also see in the VNIR range which is 350 to 1000 nanometer they have 512 element NIR-enhanced silicon silicon array, and then SWIR 1 they are calling it SWIR 1 from 1000 shortwave infrared 1000 to 1800 nanometer detector they have the again indium, gallium, arsenide detector and then SWIR 2 from 1800 nanometers to 2500 nanometer.

There also the indium, gallium, arsenide detector. So, you can see that different types of spectral features and their specifications are given. So, different types of FOV options fiber mount options are given we are going to discuss all these, wavelength accuracy is also given. So, you can see here in case a PSR 3500 wavelength accuracy is plus minus 0.5 bandwidth 0.5 of bandwidth and in case of FieldSpec, the 0.5 nanometer average, so, we can see that plus minus 1 nanometer for any line so, you can see 0.5 nanometer average error of wavelength calibration fit.

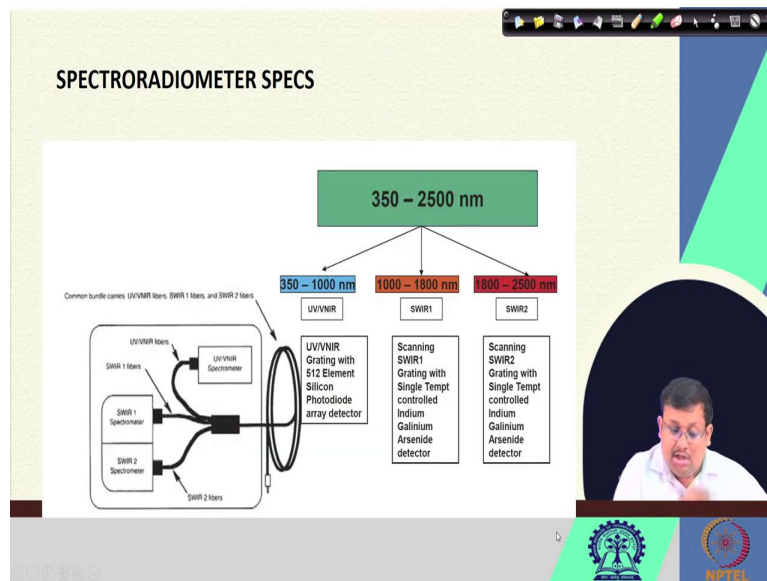
So, this is how you compare the minimum scanning time you can see here 100 milliseconds and here it is given 200 millisecond, but it is a scanning time. So, this is how, you can compare the different commercially available spectroradiometers and what are the meaning of these we are going to discuss what is spectral resolution, what is full width half maximum, we are going to discuss but depending on your objective, depending on your purpose, you can select which one serve your purpose best.

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So, if you see the spectroradiometer specs, the spectroradiometer is basically composed of these 4 or 5 components, one is a light source and then a dispersive unit or monochromator and then a detector and then there are fibers and then absorbance or reflectance standard. So, using these 5 components, we can measure the spectroradiometer generally captures the spectral information.

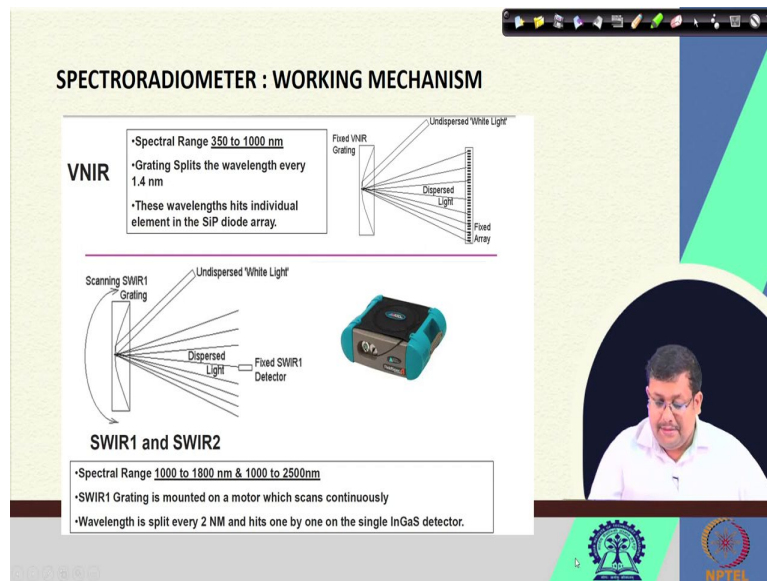
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So, if we consider this modular concept, there are 3 zones, this is actually given by the ASD FieldSpec. So, you can see they are naming it the 350 to 1000 nanometer they are naming it UV-VNIR and then 1000 to 1800 nanometer SWIR 1 and then 1800 to 2500 nanometer they are calling SWIR 2. So, as I have told you that in these UV-VNIR range they are using 512 silicon photodiode array detector. SWIR 1 they are using the indium, gallium, arsenide detector and in the SWIR 2 also they are using the indium, gallium, arsenide detector.

So, basically once the capture the common there is common bundle carriers of these a fiber optic cable, so, which capture the reflected energy and then it distributes the energy into 3 of these spectrometer, one is UV-VNIR, then SWIR 1, SWIR 2 and then they capture individually this individual region and then combine the results combined reflectance values.

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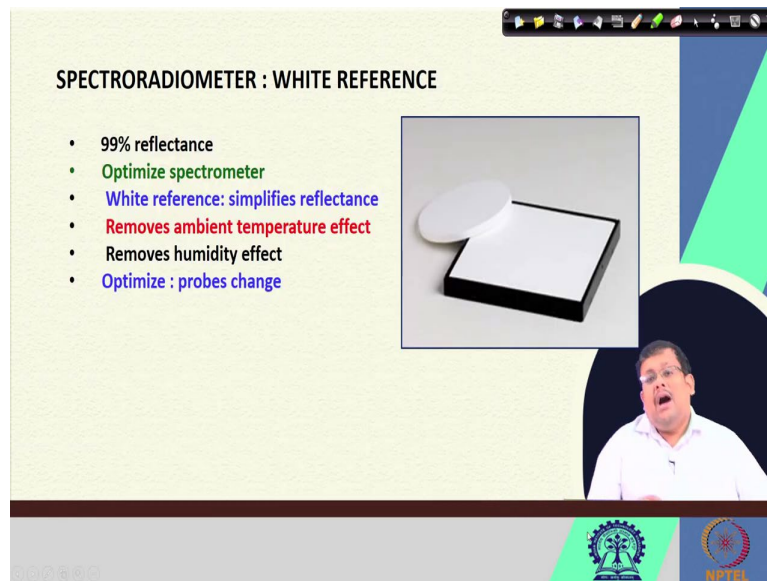


So, if you see the working mechanism, the VNIR, which is having the 512 silicon photodiode array, which is a which covers the spectral range of 350 to 1000 nanometer, here this is the fixed array and also here you can see these are undisbursed white light and fixed VNIR grating, so, undisbursed white light is divided into dispersed and then they go through this fixed array. So, grating splits the wavelength every 1.4 nanometer and these wavelengths heats individual elements of the silicon photodiode array.

When we consider the indium, gallium, arsenide detector of SWIR 1 and SWIR 2 range these they capture the spectral range of 1000 to 1800 nanometer and then 1000 to 2500 nanometer, it should be 1800 to 2500 nanometers please correct it, sorry for this so, 1800 to 2500 nanometer. So, these grating is basically mounted on a motor which scans continuously and wavelengths.

So, here the undispersed light and then the scanning SWIR 1 grating which is movable which is mounted on a motor and then there is a fixed SWIR 1 detector. So, individual dispersed lies will go through the SWIR 1 detector and get detected. So, this is how this is working mechanism of different detectors.

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**SPECTRORADIOMETER : WHITE REFERENCE**

- 99% reflectance
- Optimize spectrometer
- White reference: simplifies reflectance
- Removes ambient temperature effect
- Removes humidity effect
- Optimize : probes change

The slide features a yellow background with a blue and green geometric design on the right. An inset image shows a white circular disc and a white rectangular panel. A small video window in the bottom right corner shows a man speaking. Logos for IIT Bombay and NPTEL are visible at the bottom.

We need for any spectral measurement we need a white reference panel we call them spectral on panel. Generally, we assume that it has 99 percent reflectance. It is used for optimizing the spectrophotometer. Also, it requires for white reference for simplifying the reflectance. It helps in the removal of ambient temperature effect and also it removes the humidity effect and optimize that and when we change the probe, there are different types of probe.

When we change the probe also in between that also to standardize the instrument we require this white reference panel which is generally made up of some PTFE or poly tetra fluoro ethane, which is and also it is all it almost pure white and we assume that this panel has 99 percent reflectance.

So, we, this panel should be taken should be maintained very, very carefully, so, that there is no scratch or no soil of this white reference panels because this is the one which will be used for optimizing the spectroradiometer.



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**SPECTRAL RESOLUTION**

- Spectral sampling interval is the spacing between sample points in the spectrum.
- Sampling is independent of resolution and in ASD spectroradiometers is between 2 and 5 times per FWHM.
- The sampling interval here is 1.4 nm for the region 350 - 1000 nm and 2 nm for the region 1000 - 2500 nm.
- Spectral resolution is defined as the full-width-half-maximum (FWHM) of the instrument response to a monochromatic source.
- These spectral resolution values have been measured by calculating the FWHM of a near monochromatic peak in a spectrum acquired when viewing the output of a monochromator with the FieldSpec3<sup>®</sup> spectroradiometer.

The graph shows a bell-shaped curve representing a spectral response. The x-axis is labeled 'Wavelength' and the y-axis is 'Spectrometer Offset Corrected Raw DNS'. A vertical line marks the peak. A horizontal line at 50% of the peak height intersects the curve at two points, and the distance between these points is labeled 'Full width at half maximum (FWHM)'. A horizontal double-headed arrow above the curve indicates the 'Sampling Interval'. A vertical double-headed arrow between the peak and the 50% height line is labeled 'Spectral Resolution'.

NPTEL

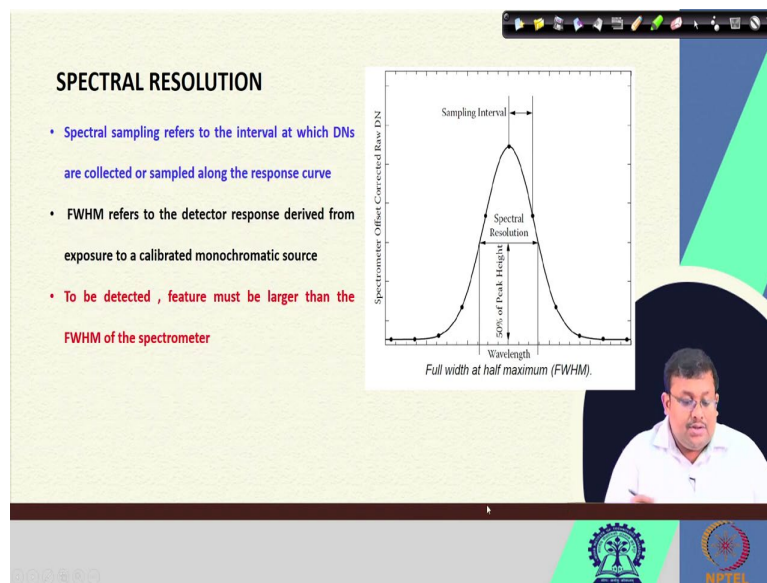
So, let us discuss what is spectral resolution? So, before we discuss the spectral resolution, let us discuss the what is spectral interval or spectral sampling interval also. So, if you see this is a spectral response from a monochromatic source and you can see here this is an absorption, absorbance spectra. So, sampling interval, so, if you consider the difference between these 2 points, the sampling interval is the spacing between sample points in the spectra. So, these are the sample points. So, the spacing is known as the sampling interval.

So, remember that the sampling is independent of the resolution in case of as far as the ASD spectroradiometers it is between 2 to 5 times per FWHM, what FWHM? FWHM means if full width half maximum. So, this is the maximum height and full width half of maximum. So, this is the 50 percent of the peak height. So, this is the peak height and this is the 50 percent of the peak height. So, this region is known as the full width at this half of maximum height or peak height. So, this is known as also the spectral resolution.

So, this is the sampling and sampling interval is generally is between 2 to 5 times per FWHM. So, within one FWHM there are 2 to 5 sampling interval. So, the sampling interval here is 1.4 nanometer for the region of 300 in case of ASD and also in case of PSR we can see the sampling interval is around 1.4 to 2 nanometer in the region between 350 to 1000 nanometer and 2 nanometer for the region of 1000 to 2500 nanometer. So, the spectral resolution so, sampling interval we can see here, what is spectral resolution? The spectral resolution is this full width half maximum.

So, spectral resolution is defined as the full width half maximum of the instrument response to a monochromatic source. This is a monochromatic source instrument response at full width half maximum. So, these spectral resolution values have been measured by calculating the full width half maximum of a near monochromatic peak in a spectrum accurate when viewing the output of a monochromator with the FieldSpec spectroradiometer. So, higher the spectral resolution that means, more narrow this full width half maximum. So, narrower this full width half maximum means higher the spectral resolution.

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So, spectral sampling generally or spectral sampling spectral interval generally refers to the interval at which the data numbers are collected or sampled along the response curve. So, there is initially this spectroradiometer captures the data number and then they convert it to the reflectance and absorbance. So, this spectral sampling refers to the interval at which the data numbers or the DNs, DNs are collected or sample along the response curve. And this FWHM refers to the detected response derived from the exposure to a calibrated monochromatic source and to be detected feature must be larger than the FWHM spectroradiometer.

So, that is why this resolution of the spectroradiometer should be, the spectral resolution of the spectroradiometer should be less than that of the spectral features which are in case of soil generally less than 20 nanometers.

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### PRE AND POST-DISPERSIVE SPECTROMETER

- **Pre-dispersive spectrometer:**
  - The ambient stray light signal can represent a large fraction of the total light signal measured by the detector
  - Therefore, it can be a major source of error
- **Post-Dispersive spectrometer**
  - The ambient stray light scattered from the sample is also collected, but with the post-dispersive instrument only ambient stray light of the same wavelength as that being measured by the detector is added to the signal.
  - Thus, the stray light signal represents a much smaller fraction of the total light signal measured by the detector resulting in an increase in instrument baseline stability

So, let us discuss this pre-dispersive and post-dispersive spectrometer in our next lecture, but I hope that in this lecture, you have gathered enough information regarding the specs of the spectroradiometer the spec, I hope that the important features of the spectroradiometer and spectral resolution, sampling interval we have already covered. And you now have an idea of how this diffuse reflectance spectrophotometer generally works.

So, in the next lecture, we are going to discuss some of the other important features and specification like pre-dispersive and post-dispersive spectrometers. We are also going to discuss the field of view and also we are going to discuss the different spectral processing. So, thank you guys.

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### REFERENCES

- [https://fsf.nerc.ac.uk/instruments/asd\\_fieldspec.shtml](https://fsf.nerc.ac.uk/instruments/asd_fieldspec.shtml)
- Viscarra-Rossel RA, McBratney AB (2008) Diffuse reflectance spectroscopy as a tool for digital soil mapping. In: Digital soil mapping with limited data. Springer, Berlin, pp 165–172
- Wadoux, A.M.J.C., Malone, B., Minasny, B., Fajardo, M., McBratney, A.B., (2021). Soil Spectral Inference with R Analysing Digital Soil Spectra using the R Programming Environment. Springer.

These are the references which I have used in these lectures, these Viscarra Rossel and McBratney. And also I would recommend you to go through these books written by Wadoux, Malone, Minasny, Fajardo, McBratney, this is a one of the recent and most comprehensive books, where you will learn how to deal with the spectral data using R. So, R is a very good free software and you should know you should utilize this software for doing a lot of different types of spectral modeling or spectro transfer modeling.

And you can do a lot of statistical analysis using R which is an open source. So, you should use this R for doing all the spectral transfer functions. And you can go with, you can see this book I am highly recommending this book that is Soil Spectral Inference with R Analyzing Digital Soil Spectra using R Programming Environment written by these authors who are known as the topmost scientist in the world as far as the soil spectroscopy is concerned.

So, you can gather more knowledge from this book if you are interested. So, thank you guys, let us meet in our next lecture, where we will discuss a different types of spectral specs of the spectroradiometers and also we are also going to discuss the spectral preprocessing. Thank you.