

# Package ‘lsatTS’

March 4, 2022

**Title** An R package to facilitate retrieval, cleaning, cross-calibration, and phenological modeling of Landsat time-series data

**Version** 0.8.0

**Description** This software package facilitates sample-based time series analysis of surface reflectance and spectral indices derived from Landsat sensors. The package includes functions that enable the extraction of the full Landsat record for point sample locations or small study regions using the Google Earth Engine accessed directly from R. Moreover, the package includes functions for (1) rigorous data cleaning, (2) cross-sensor calibration with machine learning, (3) phenological modeling, and (4) time series analysis.

**License** Modified MIT

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## R topics documented:

lsat_calc_spec_index . . . . .	2
lsat_calc_trend . . . . .	2
lsat_calibrate_rf . . . . .	3
lsat_clean_data . . . . .	5
lsat_evaluate_phenological_max . . . . .	6
lsat_export_ts . . . . .	7
lsat_fit_phenological_curves . . . . .	9
lsat_general_prep . . . . .	11
lsat_get_pixel_centers . . . . .	11
lsat_neighborhood_mean . . . . .	13
lsat_summarize_data_avail . . . . .	14
lsat_summarize_growing_seasons . . . . .	15

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lsat\_calc\_spec\_index     *Calculate Spectral Indices*


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### Description

This function computes some widely used spectral vegetation indices. Only one index can be computed at a time. Current indices include the: Normalized Difference Vegetation Index (NDVI; Rouse et al. 1974), kernel NDVI (kNDVI; Camp-Valls et al. 2020), Green NDVI (gNDVI; Gitelson and Merzlyak 1998), Soil Adjusted Vegetation Index (SAVI; Huete 1998), Wide Dynamic Range Vegetation Index (WDRVI; Gitelson 2004), Enhanced Vegetation Index (EVI; Huete et al. 2002), 2-band EVI (EVI2; Jiang et al. 2008), Near Infrared Vegetation Index (NIRv; Badgley et al. 2017), Moisture Stress Index (MSI; Rock et al. 1986), Normalized Difference Water Index (NDWI; McFeeters 1996), Normalized Difference Moisture Index (NDMI; Gao 1996), Normalized Burn Ratio (NBR, Key and Benson 1999), Normalized Difference Infrared Index (NDII; Hardisky et al. 1983), Plant Senescence Reflectance Index (PSRI; Merzlyak et al. 1999), and the Soil-Adjusted Total Vegetation Index (SATVI; Marsett et al. 2006).

### Usage

```
lsat_calc_spec_index(dt, si)
```

### Arguments

dt	Data.table containing surface reflectance data
si	Character string specifying abbreviation of the desired spectral index

### Value

The input data.table with an appended column containing the spectral index

### Examples

```
# my.dt <- lsat_calc_spec_index(my.dt, 'ndvi')
```

---

lsat\_calc\_trend     *Calculate non-parametric vegetation greenness trends*


---

### Description

This function computes a temporal trend in annual time series of vegetation greenness for each sampling site over a user-specified time period. This is a wrapper for the `zyp.yuepilon()` function from the `zyp` package. This function will iteratively pre-whiten a time series (i.e., remove temporal autocorrelation) and then compute Mann-Kendall trend tests and Theil-Sen slope indicators.

**Usage**

```
lsat_calc_trend(
  dt,
  si,
  yrs,
  yr.tolerance = 1,
  nyr.min.frac = 0.66,
  sig = 0.1,
  legend.position = c(0.8, 0.2),
  legend.dir = "horizontal"
)
```

**Arguments**

<code>dt</code>	Data.table with columns including site, year, and the vegetation index of interest
<code>si</code>	Spectral index (e.g., NDVI) for which to assess trend
<code>yrs</code>	A sequence of years (time period) over which to assess trends (e.g., 2000:2020)
<code>yr.tolerance</code>	The number of years that a site's first/last years of observations can differ from the start/end of the user-specified time period ('yrs') for a trend to be computed
<code>nyr.min.frac</code>	Fraction of years within the time period for which observations must be available if a trend is to be computed
<code>sig</code>	A p-value significance cutoff used to categories trends (e.g., 0.10)

**Value**

Data.table with summary of temporal trends by site and a histogram showing summarizing relative changes in vegetation greenness

**Examples**

```
# Forthcoming...
```

---

lsat_calibrate_rf	<i>Cross-Calibrate Landsat Sensors using Random Forests Models</i>
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**Description**

There are systematic differences in spectral indices (e.g., NDVI) among Landsat 5, 7, and 8 (Landsat Collection 1). It is important to address these differences before assessing temporal trends in spectral data. Failure to address these differences can, for instance, introduce artificial positive trends into NDVI time-series that are based on measurements from multiple Landsat sensors (Ju and Masek 2016, Roy et al. 2016, Berner et al. 2020).

This function cross-calibrates individual bands or spectral indices from Landsat 5/8 to match Landsat 7. Landsat 7 is used as a benchmark because it temporally overlaps with the other two sensors. Cross-calibration can only be performed on one band or spectral index at a time. The approach involves determining the typical reflectance at a sample during a portion of the growing season using Landsat 7 and Landsat 5/8 data that were collected the same years. A random forest model is then trained to predict Landsat 7 reflectance from Landsat 5/8 reflectance. To account for potential

seasonal and regional differences between sensors, the random forest models also include as covariates the midpoint of each 15-day period (day of year) and the spatial coordinates of each sampling sample. This approach can handle non-linear relationships, but is most suitable when working with data from 100s to preferably 1000s of sampling samples.

The specific steps to cross-calibrating sensors include: (1) Identify the years when both Landsat 7 and Landsat 5/8 measured surface reflectance at a sampling sample. (2) Pool the reflectance measurements across those years and compute 15-day moving median reflectance over the course of the growing season for each sensor and sampling sample. (3) Exclude 15-day periods with fewer than a specified number of measurements from both sets of sensors and then randomly select one remaining 15-day period from each sampling sample. (4) Split the data into sets for model training and evaluation. (5) Train Random Forest models that predict Landsat 7 reflectance based on Landsat 5/8 reflectance. The models also account for potential seasonal and regional differences between sensors by including as covariates the midpoint of each 15-day period (day of year) and the spatial coordinates of each sampling sample. (6) Fit the random forest models using the ranger package. See Berner et al. (2020) for a full description of the approach.

### Usage

```
lsat_calibrate_rf(
  dt,
  band.or.si,
  doy.rng = 152:243,
  min.obs = 5,
  train.with.highlat.data = F,
  add.predictors = NULL,
  frac.train = 0.75,
  trim = T,
  overwrite.col = F,
  outfile.id = band.or.si,
  outdir
)
```

### Arguments

<code>dt</code>	Data.table containing the band or spectral index to cross calibrate.
<code>band.or.si</code>	Character string matching the column name of the band or spectral index to cross-calibrate.
<code>doy.rng</code>	Sequence of numbers specifying the Days of Year (Julian Days) to use for model development.
<code>min.obs</code>	Minimum number of paired, seasonally-matched observations from Landsat 7 and Landsat 5/8 required to include a sampling sample.
<code>train.with.highlat.data</code>	Should the RF models be trained using an internal high-latitude dataset that sampled the Arctic and Boreal biomes?
<code>add.predictors</code>	Vector of additional predictors to use in the Random Forest models. These should be time-invariant and match column names.
<code>frac.train</code>	Fraction of samples to use for training the random forest models. The remaining samples are used for model cross-validation.
<code>trim</code>	(True/False) If true, then the percent difference between satellites is determined for each sample and then the lowest 2.5 and highest 97.5 percentiles are trimmed.

overwrite.col	(True/False) Overwrite existing column or (by default) append cross-calibrated data as a new column?
outfile.id	Identifier used when naming output files. Defaults to the input band, but can be specified if needed such as when performing Monte Carlo simulations.
outdir	Output directory (created if necessary) to which multiple files will be written. The files include the (1) fitted random forest models as R objects, (2) evaluation data in a csv file, (3) summary of model cross-validation in a csv file, and (4) multi-panel scatter plot comparing sensors pre- and post-calibration in jpeg format. If cross-calibrating both Landsat 5 and 8, then the function returns files for both sensors.

### Value

The input data.table with an appended column titled band.xcal, where "band" is your specified band or spectral index

### Examples

```
# lsat.dt <- lsat_xcal_rf(lsat.dt, band = 'ndvi', doy.rng = 152:243, min.obs = 5, frac.train = 0.75, outfile.id
```

---

lsat_clean_data	<i>Clean Landsat Data</i>
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---

### Description

This function filters out observations that exhibit: (1) clouds and optionally water and snow; (2) impossibly high reflectance ( $>1.0$ ) and abnormally low reflectance ( $<0.005$ ); (3) scene cloud cover above a user-defined threshold; (4) geometric uncertainty exceeding a user-defined threshold. More information about Landsat Quality Bands here <https://landsat.usgs.gov/collectionqualityband>

### Usage

```
lsat_clean_data(
  dt,
  cloud.max = 80,
  geom.max = 30,
  sza.max = 60,
  filter.cfmask.snow = T,
  filter.cfmask.water = T,
  filter.jrc.water = T
)
```

### Arguments

dt	Data.table generated by calling lsat_general_prep().
cloud.max	Maximum allowable cloud cover in Landsat scene (percentage).
geom.max	Maximum allowable geometric uncertainty (meters).
sza.max	Maximum allowable solar zenith angle (degrees).
filter.cfmask.snow	(TRUE/FALSE) Remove observations with CFmask flag = snow.

```
filter.cfmask.water
```

(TRUE/FALSE) Remove observations with CFmask flag = water.

```
filter.jrc.water
```

(TRUE/FALSE) Remove observations that were ever inundated based on JRC Global Surface Water Dataset.

### Value

A data.table that includes Landsat observations that met the quality control criteria.

### Examples

```
# lsat.dt <- lsat_clean_dt(lsat.dt, cloud.max=80, geom.max=30, sza.max=60, filter.snow = T, filter.water = T)
```

---

```
lsat_evaluate_phenological_max
```

*Evaluate Estimates of Annual Phenological Maximum*

---

### Description

Assess how the number of annual Landsat observations impacts estimates of annual maximum vegetation greenness derived from raw observations and phenological modeling. The algorithm computes annual maximum vegetation greenness using site x years with a user-specific number of observations and then compares these with estimates derived when using progressively smaller subsets of observations. This lets you determine the degree to which annual estimates of maximum vegetation greenness are impacted by the number of available observations.

### Usage

```
lsat_evaluate_phenological_max(
  dt,
  si,
  min.frac.of.max = 0.75,
  zscore.thresh = 3,
  min.obs = 6,
  reps = 10,
  outdir = NA
)
```

### Arguments

dt	Data.table output from lsat_fit_phenologic_curves().
si	Character string specifying the spectral index (e.g., NDVI) to evaluate.
min.frac.of.max	Numeric threshold (0-1) that defines the "growing season" as the seasonal window when the phenological curves indicate the VI is within a specified fraction of the maximum VI. In other words, an observation is considered to be from the "growing season" when the VI is within a user-specified fraction of the curve-fit growing season maximum VI.
zscore.thresh	Numeric threshold specifying the Z-score value beyond which individual observations are filtered before computing the maximum VI.

min.obs	Minimum number of observations needed for a site x year to be included in the evaluation (Default = 10)
reps	Number of times to bootstrap the assessment (Default = 10)
outdir	If desired, specify the output directory where evaluation data and figure should be written. If left as NA, then no output is only displayed in the console and not written to disk.

### Value

Data.table

### Examples

```
# Forthcoming...
```

---

lsat_export_ts	<i>Export Surface Reflectance Time-Series from the Landsat C2 record using rgee</i>
----------------	---

---

### Description

This function exports the surface reflectance time series for a set of point-coordinates from the whole Landsat C2 record using the Google Earth Engine (account required). These resulting time-series can then be processed using the remainder of the lsatTS package workflow. For polygon geometries consider using `lsat_get_pixel_centers()` to generate pixel center coordinates for all pixels within a given polygon first.

### Usage

```
lsat_export_ts(
  pixel_coords_sf,
  sample_id_from = "sample_id",
  chunks_from = NULL,
  this_chunk_only = NULL,
  max_chunk_size = 250,
  drive_export_dir = "lsatTS_export",
  file_prefix = "lsatTS_export",
  startJulian = 152,
  endJulian = 243,
  start_date = "1984-01-01",
  end_date = "today",
  BUFFER_DIST = 0,
  SCALE = 30,
  MASK_VALUE = 0
)
```

## Arguments

pixel_coords_sf	Simple feature object of point coordinates for the sample.
sample_id_from	The column name that specifies the unique sample identifier in pixel_coords_sf (defaults to "sample_id" as generated by lsat_get_pixel_centres).
chunks_from	Column name in pixel_coords_sf to divide the exports into chunks. Over-rides chunk division by size (see max_chunk_size).
this_chunk_only	Name of a specific chunk to be exported. Useful for re-exporting a single chunk should the export fail for some reason.
max_chunk_size	Maximum number of sample coordinates to be exported in each chunk. Defaults to 250.
drive_export_dir	Folder on the user's Google Drive to export the records too. Defaults to "lsatTS_export".
file_prefix	Optional file_prefix for the exported files.
startJulian	Optional first day of year to extract for. Defaults to 152.
endJulian	Optional Last day of year to extract for. Defaults to 243.
start_date	Optional extraction start date (as string, format YYYY-MM-DD). Defaults to "1984-01-01".
end_date	Optional extraction end date (as string, format YYYY-MM-DD). Defaults to today's date.
BUFFER_DIST	Buffer distance around sample coordinates. Uses lsat_get_pixel_centers() to find all Landsat pixel centers around each point in pixel_coords_sf - within the distance specified here (square buffer). Can be slow for large number of points. Defaults to 0 m.
SCALE	Scale for extraction. Defaults to 30 m nominal Landsat pixel size).
MASK_VALUE	Optional masking value for global surface water mask. Defaults to 0.

## Details

Please note: Unlike the other functions in this package, this function does NOT return the time-series as an object, instead it returns a list of the tasks issued for the export. The actual time-series are exported as CSV objects by the EE to the user's Google Drive. This way of exporting allows for a more efficient scheduling, larger exports, and does not require the R session to continue to run in the background while the requests are processed on the EE.

The progress of the exports can be monitored using the list of tasks returned by this function in combination with the ee\_monitoring() function from rgee, or using the task overview in the web code editor of the EE (<https://code.earthengine.google.com>).

## Value

List of rgee tasks.

## Author(s)

Jakob J. Assmann and Richard Massey



## Examples

```
# Using sf, dplyr and rgee
library(sf)
library(dplyr)
library(rgee)

# Initialize EE
ee_initialize()

# Generate test points
test_points_sf <- st_sfc(st_point(c(-149.6026, 68.62574)),
                        st_point(c(-149.6003, 68.62524)),
                        st_point(c(-75.78057, 78.87038)),
                        st_point(c(-75.77098, 78.87256)),
                        st_point(c(-20.56182, 74.47670)),
                        st_point(c(-20.55376, 74.47749)), crs = 4326) %>%

  st_sf() %>%
  mutate(sample_id = c("toolik_1",
                        "toolik_2",
                        "ellesmere_1",
                        "ellesmere_1",
                        "zackenberg_1",
                        "zackenberg_2"),
         region = c("toolik", "toolik",
                    "ellesmere", "ellesmere",
                    "zackenberg", "zackenberg"))

# Export time-series using lsat_export_ts()
task_list <- lsat_export_ts(test_points_sf)

# Export time-series using with a chunk size of 2
task_list <- lsat_export_ts(test_points_sf, max_chunk_size = 2)

# Export time-series in chunks by column
task_list <- lsat_export_ts(test_points_sf, chunks_from = "region")
```

---

lsat\_fit\_phenological\_curves

*Characterize Land Surface Phenology using Vegetation Index Time Series*

---

## Description

This function characterizes seasonal land surface phenology at each sample site using time series of spectral vegetation indices (e.g., NDVI). The underlying algorithm was constructed to facilitate estimating annual maximum vegetation greenness (spectral index). The function returns information about typical phenology at a sample site and about the timing of an individual observation relative. Please note that this function was designed for situations where the seasonal phenology is hump shaped. If you are using a spectral index that is typically negative (e.g., Normalized Difference Water Index) then multiply the index by -1 before running this function, then back-transform your index after running the `lsat_summarize_growing_seasons()` function.

**Usage**

```
lsat_fit_phenological_curves(
  dt,
  si,
  window.yrs = 11,
  window.min.obs = 20,
  si.min = 0.15,
  spar = 0.73,
  pcnt.dif.thresh = 30,
  weight = T,
  spl.fit.outfile = F,
  progress = T,
  test.run = F
)
```

**Arguments**

<code>dt</code>	Data.table with a multi-year time series a vegetation index
<code>si</code>	Character string specifying the spectral index (e.g., NDVI) to use for determining surface phenology. This must correspond to an existing column in the data.table.
<code>window.yrs</code>	Number specifying the focal window width in years that is used when pooling data to fit cubic splines (use odd numbers).
<code>window.min.obs</code>	Minimum number of focal window observations necessary to fit a cubic spline.
<code>si.min</code>	Minimum value of spectral index necessary for observation to be used when fitting cubic splines. Defaults to 0.15 which for NDVI is about when plants are present.
<code>spar</code>	Smoothing parameter passed to <code>smooth.spline()</code> , typically around 0.65 - 0.75 for this application.
<code>pcnt.dif.thresh</code>	Allowable percent difference (0-100) between individual observations and fitted cubic spline. Observations that differ by more than this threshold are filtered out and the cubic spline is iteratively refit.
<code>weight</code>	When fitting the cubic splines, should individual observations be weighted by their year of acquisition relative to the focal year? If so, each observation is weighted by $\exp(-0.25 * n.yrs.from.focal)$ when fitting the cubic splines.
<code>spl.fit.outfile</code>	(Optional) Name of output csv file containing the fitted cubic splines for each sample site. Useful for subsequent visualization
<code>progress</code>	(TRUE/FALSE) Print a progress report?
<code>test.run</code>	(TRUE/FALSE) If TRUE, then algorithm is run using a small random subset of data and only a figure is output. This is used for model parameterization.

**Value**

Data.table that provides, for each observation, information on the phenological conditions for that specific day of year during the focal period. These data can then be used to estimate annual maximum spectral index and other growing season metrics using `lsat_summarize_growing_season()`. A figure is also generated that shows observation points and phenological curves for nine random sample locations.

**Examples**

```
# To come...
```

---

lsat_general_prep	<i>Prepare Landsat Data</i>
-------------------	-----------------------------

---

**Description**

This function parses sample site coordinates and time period of each measurement, scales band values, and formats column names as needed for subsequent analysis using the lsatTS package.

**Usage**

```
lsat_general_prep(dt)
```

**Arguments**

dt                      Data.table with Landsat data exported from Google Earth Engine using lsat\_export\_ts()

**Value**

Data.table with formatted and scaled values.

**Examples**

```
# parsed.dt <- lsat_general_prep(gee.dt)
```

---

lsat_get_pixel_centers	<i>Get Landsat 8 pixels centers within a polygon or around a point coordinate with a buffer</i>
------------------------	---

---

**Description**

A convenience helper function that determines the Landsat 8 grid (pixel) centers within a polygon and an optional buffer. It can also be applied to a single point to retrieve all pixels within a surrounding buffer.

**Usage**

```
lsat_get_pixel_centers(
  polygon_sf,
  pixel_prefix = "pixel",
  pixel_prefix_from = NULL,
  buffer = 15,
  plot_map = F,
  lsat_WRS2_scene_bounds = NULL
)
```

**Arguments**

<code>polygon_sf</code>	Simple feature with a simple feature collection of type "sfc_POLYGON" containing a single polygon geometry. Alternatively, a simple feature containing a simple feature collection of type 'sfc_POINT' with a single point.
<code>pixel_prefix</code>	Prefix for the generated pixel ids. Defaults to "pixel".
<code>pixel_prefix_from</code>	Optional, column name in simple feature to specify pixel_prefix. Overrides "pixel_prefix" argument.
<code>buffer</code>	Buffer surrounding the geometry to be included. Specified in m. Defaults to 15 m, the nominal Landsat pixel size.
<code>plot_map</code>	Optional, default is FALSE. If TRUE the retrieved pixel centers and the polygon are plotted on a summer Landsat 8 image (grey-scale red band) using mapview. If a character is supplied an additional output to a file is generated (png, pdf, and jpg supported, see mapview::mapshot). Note: Both slow down the execution of this function dramatically, especially for large polygons. Only useful in interactive sessions.
<code>lsat_WRS2_scene_bounds</code>	File path to the Landsat WRS2 path row scene boundaries. If not specified the boundaries are downloaded to a temporary file when the function is executed the first time during a session. To avoid future downloads, the file may be downloaded manually and it's file path specified using this argument. The file can be found here: <a href="https://prd-wret.s3.us-west-2.amazonaws.com/assets/palladium/production/atoms/files/WRS2_bound_world_0.kml">https://prd-wret.s3.us-west-2.amazonaws.com/assets/palladium/production/atoms/files/WRS2_bound_world_0.kml</a> See also: <a href="https://www.usgs.gov/core-science-systems/nli/landsat/landsat-shapefiles-and-kml-files">https://www.usgs.gov/core-science-systems/nli/landsat/landsat-shapefiles-and-kml-files</a>

**Details**

Does not work for large polygons. The default maximum number of pixels set by the GEE is 10000000. Consider whether extraction for a large polygon is a good idea, if yes split the polygon into manageable chunks.

For the unlikely case that a polygon exceeds the boundaries of the Landsat tile closest to the polygon's center, the polygon is clipped at the boundaries of the Landsat tile and a warning is issued. Again, if this is the case, consider processing smaller polygons instead.

Please note that the approximation of tile overlap with polygon generates a warning by sf that the coordinates are assumed to be planar. This can be ignored.

**Value**

sfc of point geometries for Landsat 8 pixel centers within the polygon or the buffer around the point coordinate. For use in `lsat_export_ts()`.

**Author(s)**

Jakob J. Assmann

**Examples**

```
# Using sf, dplyr, rgee and purr
library(sf)
library(dplyr)
library(rgee)
```

```

library(purrr)

# Initialize EE
ee_initialize()

# Specify a region to retrieve pixel centers for
test_poly_sf <- list(matrix(c(-138.90125, 69.58413,
                             -138.88988, 69.58358,
                             -138.89147, 69.58095,
                             -138.90298, 69.57986,
                             -138.90125, 69.58413),
                             ncol = 2, byrow = TRUE)) %>%
  st_polygon() %>%
  st_sfc(crs = 4326) %>%
  st_sf()

# Retrieve pixel centers and plot to mapview
pixels <- lsat_get_pixel_centers(test_poly_sf, plot_map = TRUE)

## Ge pixel centers for multiple regions
# Create multi-polygon sf
ellesmere <- st_polygon(list(matrix(c(-75.78526, 78.86973,
                                     -75.78526, 78.87246,
                                     -75.77116, 78.87246,
                                     -75.77116, 78.86973,
                                     -75.78526, 78.86973),
                                     ncol = 2, byrow = TRUE)))
zackenberg <- st_polygon(list(matrix(c(-20.56254, 74.47469,
                                       -20.56254, 74.47740,
                                       -20.55242, 74.47740,
                                       -20.55242, 74.47469,
                                       -20.56254, 74.47469),
                                       ncol = 2, byrow = TRUE)))
toolik <- st_polygon(list(matrix(c(-149.60686, 68.62364,
                                   -149.60686, 68.62644,
                                   -149.59918, 68.62644,
                                   -149.59918, 68.62364,
                                   -149.60686, 68.62364),
                                   ncol = 2, byrow = TRUE)))
test_regions_sf <- st_sfc(ellesmere, zackenberg, toolik, crs = 4326) %>%
  st_sf() %>%
  mutate(region = c("ellesmere", "zackenberg", "toolik"))

# Split and map lsat_get_pixel_centers using dplyr and purrr
pixel_list <- test_regions_sf %>%
  split(.$region) %>%
  map(lsat_get_pixel_centers,
      pixel_prefix_from = "region") %>%
  bind_rows()

```

---

lsat\_neighborhood\_mean

*Compute Neighborhood Average Landsat Surface Reflectance*


---

**Description**

For each band, this function computes average surface reflectance across neighboring voxels at a sample site. Use this function when working with Landsat data extracted for buffered points. Also, make sure to have previously cleaning the individual observations using `lsat_clean_data()`.

**Usage**

```
lsat_neighborhood_mean(dt)
```

**Arguments**

`dt` A data.table containing coincident surface reflectance measurements for multiple Landsat pixels at each sample site.

**Value**

A data.table with average surface reflectance

**Examples**

```
# ngb.avg.dt <- lsat_ngb_mean(my.dt)
```

---

```
lsat_summarize_data_avail
```

*Summarize Availability of Landsat for each Sample Site*

---

**Description**

This little function summarizes the temporal period and availability of observations at each sample site

**Usage**

```
lsat_summarize_data_avail(dt)
```

**Arguments**

`dt` Data.table with columns named "sample.id" and "year"

**Value**

Data.table summarizing for each site the first, last, and number of years with observations, the minimum and maximum number of observations in a year, and the total number of observations across years. Also returns a figure showing the median (2.5 and 97.5 percentiles) number of observations per sample site across years for each Landsat satellite.

**Examples**

```
# summary.dt <- lsat_summarize_data_avail(dt)
```

---

`lsat_summarize_growing_seasons`*Summarize a Vegetation Index for each Growing Season*

---

### Description

This function not only computes mean, median, and 90th percentile of a spectral index (SI) using observations for a user-specified "growing season," but also estimates the annual maximum VI and associated day of year using phenology modeling and growing season observations.

### Usage

```
lsat_summarize_growing_seasons(  
  dt,  
  si,  
  min.frac.of.max = 0.75,  
  zscore.thresh = 3  
)
```

### Arguments

<code>dt</code>	Data.table generated by the function <code>lsat_fit_phenologic_curves()</code> .
<code>si</code>	Character string specifying the vegetation index to summarize (e.g., NDVI).
<code>min.frac.of.max</code>	Numeric threshold (0-1) that defines the "growing season" as the seasonal window when the phenological curves indicate the SI is within a specified fraction of the maximum SI. In other words, an observation is considered to be from the "growing season" when the SI is within a user-specified fraction of the curve-fit growing season maximum SI.
<code>zscore.thresh</code>	Numeric threshold specifying the Z-score value beyond which individual observations are filtered before summarizing growing season SI.

### Value

Data.table summarizing annual growing season conditions based on a vegetation index.

### Examples

```
# Forthcoming...
```