

Package ‘epwshiftr’

July 7, 2020

Title Create Future 'EnergyPlus' Weather Files using 'CMIP6' Data

Version 0.1.0

Description Query, download climate projection data from the 'CMIP6' (Coupled Model Intercomparison Project Phase 6) project <<https://pcmdi.llnl.gov/CMIP6>> in the 'ESGF' (Earth System Grid Federation) platform <<https://esgf.llnl.gov>>, and create future 'EnergyPlus' <<https://energyplus.net>> Weather ('EPW') files adjusted from climate changes using data from Global Climate Models ('GCM').

Imports checkmate, data.table (>= 1.12.4), eplusr, future.apply, fst, jsonlite, progressr, psychrolib, rappidirs, RNetCDF, units

Suggests testthat, pingr, covr

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Encoding UTF-8

LazyData true

URL <https://github.com/ideas-lab-nus/epwshiftr>

BugReports <https://github.com/ideas-lab-nus/epwshiftr/issues>

RoxygenNote 7.0.2

Collate 'coord.R' 'utils.R' 'epwshiftr-package.R' 'esgf.R' 'morph.R'
'netcdf.R'

NeedsCompilation no

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R topics documented:

| | |
|-----------------------------|---|
| epwshiftr-package | 2 |
| esgf_query | 3 |
| extract_data | 7 |

| | |
|----------------------------|----|
| future_epw | 9 |
| get_data_dir | 10 |
| get_data_node | 11 |
| init_cmip6_index | 11 |
| load_cmip6_index | 15 |
| match_coord | 16 |
| morphing_epw | 17 |
| set_cmip6_index | 18 |
| summary_database | 19 |

Index**22**

epwshifttr-package

*epwshifttr: Create future EnergyPlus Weather files using CMIP6 data***Description**

Query, download climate change projection data from the [CMIP6 \(Coupled Model Intercomparison Project Phase 6\) project](#) in the [ESGF \(Earth System Grid Federation\) platform](#), and create future [EnergyPlus](#) Weather (EPW) files adjusted from climate changes using data from Global Climate Models (GCM).

Package options

- `epwshifttr.verbose`: If TRUE, more detailed message will be printed. Default: FALSE.
- `epwshifttr.dir`: The directory to store package data, including CMIP6 model output file index and etc. If not set, the current user data directory will be used.

Author(s)

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See Also

Useful links:

- <https://github.com/ideas-lab-nus/epwshifttr>
- Report bugs at <https://github.com/ideas-lab-nus/epwshifttr/issues>

esgf_query*Query CMIP6 data using ESGF search RESTful API*

Description

Query CMIP6 data using ESGF search RESTful API

Usage

```
esgf_query(
  activity = "ScenarioMIP",
  variable = c("tas", "tasmax", "tasmin", "hurs", "hursmax", "hursmin", "pr", "rsds",
    "rlds", "psl", "sfcWind", "clt"),
  frequency = "day",
  experiment = c("ssp126", "ssp245", "ssp370", "ssp585"),
  source = c("AWI-CM-1-1-MR", "BCC-CSM2-MR", "CESM2", "CESM2-WACCM", "EC-Earth3",
    "EC-Earth3-Veg", "GFDL-ESM4", "INM-CM4-8", "INM-CM5-0", "MPI-ESM1-2-HR",
    "MRI-ESM2-0"),
  variant = "r1i1p1f1",
  replica = FALSE,
  latest = TRUE,
  resolution = c("100 km", "50 km"),
  type = "Dataset",
  limit = 10000L,
  data_node = NULL
)
```

Arguments

| | |
|----------|---|
| activity | A character vector indicating activity identifiers. Default: "ScenarioMIP". Possible values: <ul style="list-style-type: none"> • "AerChemMIP": Aerosols and Chemistry Model Intercomparison Project, • "C4MIP": Coupled Climate Carbon Cycle Model Intercomparison Project, • "CDRMIP": Carbon Dioxide Removal Model Intercomparison Project, • "CFMIP": Cloud Feedback Model Intercomparison Project, • "CMIP": CMIP DECK: 1pctCO₂, abrupt4xCO₂, amip, esm-piControl, esm-historical, historical, and piControl experiments, • "CORDEX": Coordinated Regional Climate Downscaling Experiment, • "DAMIP": Detection and Attribution Model Intercomparison Project, • "DCPP": Decadal Climate Prediction Project, • "DynVarMIP": Dynamics and Variability Model Intercomparison Project, • "FAFMIP": Flux-Anomaly-Forced Model Intercomparison Project, • "GMMIP": Global Monsoons Model Intercomparison Project, • "GeoMIP": Geoengineering Model Intercomparison Project, • "HighResMIP": High-Resolution Model Intercomparison Project, |
|----------|---|

| | |
|-----------|---|
| | <ul style="list-style-type: none"> • "ISMIP6": Ice Sheet Model Intercomparison Project for CMIP6, • "LS3MIP": Land Surface, Snow and Soil Moisture, • "LUMIP": Land-Use Model Intercomparison Project, • "OMIP": Ocean Model Intercomparison Project, • "PAMIP": Polar Amplification Model Intercomparison Project, • "PMIP": Palaeoclimate Modelling Intercomparison Project, • "RFMIP": Radiative Forcing Model Intercomparison Project, • "SIMIP": Sea Ice Model Intercomparison Project, • "ScenarioMIP": Scenario Model Intercomparison Project, • "VIACSAB": Vulnerability, Impacts, Adaptation and Climate Services Advisory Board, • "VolMIP": Volcanic Forcings Model Intercomparison Project |
| variable | <p>A character vector indicating variable identifiers. The 12 most related variables for EPW are set as defaults. If NULL, all possible variables are returned. Default: c("tas", "tasmax", "tasmin", "hurs", "hursmax", "hursmin", "psl", "rss", "rls", "sfcWind", "pr", "rsds", "rlds", "clt"), where:</p> <ul style="list-style-type: none"> • tas: Near-surface (usually, 2 meter) air temperature, units: K. • tasmax: Maximum near-surface (usually, 2 meter) air temperature, units: K. • tasmin: Minimum near-surface (usually, 2 meter) air temperature, units: K. • hurs: Near-surface relative humidity, units: %. • hursmax: Maximum near-surface relative humidity, units: %. • hursmin: Minimum near-surface relative humidity, units: %. • psl: Sea level pressure, units: Pa. • rsds: Surface downwelling shortwave radiation, units: W m-2. • rlds: Surface downwelling longwave radiation, units: W m-2. • sfcWind: Near-surface (usually, 10 meters) wind speed, units: m s-1. • pr: Precipitation, units: kg m-2 s-1. • clt: Total cloud area fraction for the whole atmospheric column, as seen from the surface or the top of the atmosphere. Units: %. |
| frequency | <p>A character vector of sampling frequency. If NULL, all possible frequencies are returned. Default: "day". Possible values:</p> <ul style="list-style-type: none"> • "1hr": sampled hourly, • "1hrCM": monthly-mean diurnal cycle resolving each day into 1-hour means, • "1hrPt": sampled hourly, at specified time point within an hour, • "3hr": sampled every 3 hours, • "3hrPt": sampled 3 hourly, at specified time point within the time period, • "6hr": sampled every 6 hours, • "6hrPt": sampled 6 hourly, at specified time point within the time period, • "day": daily mean samples, • "dec": decadal mean samples, • "fx": fixed (time invariant) field, |

| | |
|------------|--|
| | <ul style="list-style-type: none"> • "mon": monthly mean samples, • "monC": monthly climatology computed from monthly mean samples, • "monPt": sampled monthly, at specified time point within the time period, • "subhrPt": sampled sub-hourly, at specified time point within an hour, • "yr": annual mean samples, • "yrPt": sampled yearly, at specified time point within the time period |
| experiment | A character vector indicating root experiment identifiers. The Tier-1 experiment of activity ScenarioMIP are set as defaults. If NULL, all possible experiments are returned. Default: c("ssp126", "ssp245", "ssp370", "ssp585"). |
| source | A character vector indicating model identifiers. Defaults are set to 11 sources which give outputs of all 4 experiments of activity ScenarioMIP with daily frequency, i.e. "AWI-CM-1-1-MR", "BCC-CSM2-MR", "CESM2", "CESM2-WACCM", "EC-Earth3", "EC-Earth3-Veg", "GFDL-ESM4", "INM-CM4-8", "INM-CM5-0", "MPI-ESM1-2-HR" and "MRI-ESM2-0". If NULL, all possible sources are returned. |
| variant | <p>A character vector indicating label constructed from 4 indices stored as global attributes in format r<k>i<l>p<m>f<n> described below. Default: "r1i1p1f1". If NULL, all possible variants are returned.</p> <ul style="list-style-type: none"> • r: realization_index (<k>) = realization number (integer >0) • i: initialization_index (<l>) = index for variant of initialization method (integer >0) • p: physics_index (<m>) = index for model physics variant (integer >0) • f: forcing_index (<n>) = index for variant of forcing (integer >0) |
| replica | Whether the record is the "master" copy, or a replica. Use FALSE to return only originals and TRUE to return only replicas. Default: FALSE. |
| latest | Whether the record is the latest available version, or a previous version. Use TRUE to return only the latest version of all records and FALSE to return previous versions. Default: FALSE. |
| resolution | A character vector indicating approximate horizontal resolution. Default: c("50 km", "100 km"). If NULL, all possible resolutions are returned. |
| type | A single string indicating the intrinsic type of the record. Should be either "Dataset" or "File". Default: "Dataset". |
| limit | An integer indicating the maximum of matched records to return. Should be <= 10,000. Default: 10000. |
| data_node | A character vector indicating data nodes to be queried. Default to NULL, which means all possible data nodes. |

Details

The Earth System Grid Federation (ESGF) is an international collaboration for the software that powers most global climate change research, notably assessments by the Intergovernmental Panel on Climate Change (IPCC).

The ESGF search service exposes a RESTful URL that can be used by clients to query the contents of the underlying search index, and return results matching the given constraints. ributed capabilities

of the ESGF search, the URL at any Index Node can be used to query that Node only, or all Nodes in the ESGF system. `esgf_query()` uses the [LLNL \(Lawrence Livermore National Laboratory\)](#) Index Node.

Value

A `data.table::data.table` with an attribute named `response` which is a list converted from json response. If no matched data is found, an empty `data.table` is returned. Otherwise, the columns of returned data varies based on the type:

- If "Dataset", returned columns are:

| No. | Column | Type | Description |
|-----|--------------------|-----------|--|
| 1 | dataset_id | Character | Dataset universal identifier |
| 2 | mip_era | Character | Activity's associated CMIP cycle. Will always be "CMIP6" |
| 3 | activity_drs | Character | Activity DRS (Data Reference Syntax) |
| 4 | institution_id | Character | Institution identifier |
| 5 | source_id | Character | Model identifier |
| 6 | experiment_id | Character | Root experiment identifier |
| 7 | member_id | Character | A compound construction from sub_experiment_id and variant_label |
| 8 | table_id | Character | Table identifier, i.e. sampling frequency identifier |
| 9 | frequency | Character | Sampling frequency |
| 10 | grid_label | Character | Grid identifier |
| 11 | version | Character | Approximate date of model output file |
| 12 | nominal_resolution | Character | Approximate horizontal resolution |
| 13 | variable_id | Character | Variable identifier |
| 14 | variable_long_name | Character | Variable long name |
| 15 | variable_units | Character | Units of variable |
| 16 | data_node | Character | Data node to download the model output file |
| 17 | dataset_pid | Character | A unique string that helps identify the dataset |

- If "File", returned columns are:

| No. | Column | Type | Description |
|-----|--------------------|-----------|--|
| 1 | file_id | Character | Model output file universal identifier |
| 2 | dataset_id | Character | Dataset universal identifier |
| 3 | mip_era | Character | Activity's associated CMIP cycle. Will always be "CMIP6" |
| 4 | activity_drs | Character | Activity DRS (Data Reference Syntax) |
| 5 | institution_id | Character | Institution identifier |
| 6 | source_id | Character | Model identifier |
| 7 | experiment_id | Character | Root experiment identifier |
| 8 | member_id | Character | A compound construction from sub_experiment_id and variant_label |
| 9 | table_id | Character | Table identifier, i.e. sampling frequency identifier |
| 10 | frequency | Character | Sampling frequency |
| 11 | grid_label | Character | Grid identifier |
| 12 | version | Character | Approximate date of model output file |
| 13 | nominal_resolution | Character | Approximate horizontal resolution |
| 14 | variable_id | Character | Variable identifier |

| | | | |
|----|--------------------|-----------|---|
| 15 | variable_long_name | Character | Variable long name |
| 16 | variable_units | Character | Units of variable |
| 17 | datetime_start | POSIXct | Start date and time of simulation |
| 18 | datetime_end | POSIXct | End date and time of simulation |
| 19 | file_size | Character | Model output file size in Bytes |
| 20 | data_node | Character | Data node to download the model output file |
| 21 | file_url | Character | Model output file download url from HTTP server |
| 22 | tracking_id | Character | A unique string that helps identify the output file |

References

https://github.com/ESGF/esgf.github.io/wiki/ESGF_Search_REST_API

Examples

```
## Not run:
esgf_query(variable = "rss", experiment = "ssp126", resolution = "100 km", limit = 1)

esgf_query(variable = "rss", experiment = "ssp126", type = "File", limit = 1)

## End(Not run)
```

extract_data

Extract data

Description

extract_data() takes an epw_cmip6_coord object generated using [match_coord\(\)](#) and extracts CMIP6 data using the coordinates and years of interest specified.

Usage

```
extract_data(
  coord,
  years = NULL,
  unit = FALSE,
  out_dir = NULL,
  by = NULL,
  keep = is.null(out_dir),
  compress = 100
)
```

Arguments

| | |
|----------|--|
| coord | An epw_cmip6_coord object created using match_coord() |
| years | An integer vector indicating the target years to be included in the data file. All other years will be excluded. If NULL, no subsetting on years will be performed. Default: NULL. |
| unit | If TRUE, units will be added to values using units::set_units() . |
| out_dir | The directory to save extracted data using fst::write_fst() . If NULL, all data will be kept in memory by default. Default: NULL. |
| by | A character vector of variable names used to split data during extraction. Should be a subset of: <ul style="list-style-type: none"> • "experiment": root experiment identifiers • "source": model identifiers • "variable": variable identifiers • "activity": activity identifiers • "frequency": sampling frequency • "variant": variant label • "resolution": approximate horizontal resolution If NULL and out_dir is given, file name data.fst will be used. Default: NULL. |
| keep | Whether keep extracted data in memory. Default: TRUE if out_dir is NULL, and FALSE otherwise. |
| compress | A single integer in the range 0 to 100, indicating the amount of compression to use. Lower values mean larger file sizes. Default: 100. |

Details

`extract_data()` uses [future.apply](#) underneath. You can use your preferable [future](#) backend to speed up data extraction in parallel. By default, `extract_data()` uses [future::sequential](#) backend, which runs things in sequential.

Value

An epw_cmip6_data object, which is basically a list of 3 elements:

- epw: An [eplusr::Epw](#) object whose longitude and latitute are used to extract CMIP6 data. It is the same object as created in [match_coord\(\)](#)
- meta: A list containing basic meta data of input EPW, including city, state_province, country, latitute and longitude.
- data: An empty [data.table::data.table\(\)](#) if keep is FALSE or a [data.table::data.table\(\)](#) of 12 columns if keep is TRUE:

| No. | Column | Type | Description |
|-----|----------------|-----------|--------------------------------------|
| 1 | activity_drs | Character | Activity DRS (Data Reference Syntax) |
| 2 | institution_id | Character | Institution identifier |
| 3 | source_id | Character | Model identifier |
| 4 | experiment_id | Character | Root experiment identifier |

| | | | |
|----|-------------|-----------|--|
| 5 | member_id | Character | A compound construction from sub_experiment_id and variant_label |
| 6 | table_id | Character | Table identifier |
| 7 | lat | Double | Latitude of extracted location |
| 8 | lon | Double | Latitude of extracted location |
| 9 | variable | Character | Variable identifier |
| 10 | description | Character | Variable long name |
| 11 | units | Character | Units of variable |
| 12 | value | Double | Start date and time of simulation |

Examples

```
## Not run:
coord <- match_coord("path_to_an_EPW")
extract_data(coord, years = 2030:2060)

## End(Not run)
```

future_epw

Create future EPW files using morphed data

Description

Create future EPW files using morphed data

Usage

```
future_epw(
  morphed,
  by = c("experiment", "source", "interval"),
  dir = ".",
  separate = TRUE,
  overwrite = FALSE
)
```

Arguments

| | |
|---------|--|
| morphed | An epw_cmip6_morphed object created using morphing_epw() . |
| by | A character vector of columns to be used as grouping variables when creating EPW files. Should be a subset of: <ul style="list-style-type: none"> • "experiment": root experiment identifiers • "source": model identifiers • "variable": variable identifiers • "activity": activity identifiers • "frequency": sampling frequency • "variant": variant label |

| | |
|-----------|---|
| | <ul style="list-style-type: none"> • "resolution": approximate horizontal resolution • "longitude": averaged longitude of input data • "latitude": averaged latitude of input data |
| dir | The parent directory to save the generated EPW files. If not exist, it will be created first. Default: ".", i.e., current working directory. |
| separate | If TRUE, each EPW file will be saved into a separate folder using grouping variables specified in by. |
| overwrite | If TRUE, overwrite existing files if they exist. Default: FALSE. |

Value

A list of generated [eplusr::Epw](#) objects, invisibly

| | |
|--------------|---|
| get_data_dir | <i>Get the path of directory where epwshiftr data is stored</i> |
|--------------|---|

Description

If option `epwshiftr.dir` is set, use it. Otherwise, get package data storage directory using [rappdirs::user_data_dir\(\)](#).

Usage

```
get_data_dir()
```

Value

A single string.

Examples

```
options(epwshiftr.dir = tempdir())
get_data_dir()
```

| | |
|----------------------------|--|
| <code>get_data_node</code> | <i>Get data nodes which store CMIP6 output</i> |
|----------------------------|--|

Description

Get data nodes which store CMIP6 output

Usage

```
get_data_node(speed_test = FALSE, timeout = 3)
```

Arguments

| | |
|-------------------------|---|
| <code>speed_test</code> | If TRUE, use pingr::ping() to perform connection speed test on each data node. A ping column is appended in returned data.table which stores each data node response in milliseconds. This feature needs pingr package already installed. Default: FALSE. |
| <code>timeout</code> | Timeout for a ping response in seconds. Default: 3. |

Value

A `data.table::data.table()` of 2 or 3 (when `speed_test` is TRUE) columns:

| Column | Type | Description |
|------------------------|-----------|---|
| <code>data_node</code> | character | Web address of data node |
| <code>status</code> | character | Status of data node. "UP" means OK and "DOWN" means currently not available |
| <code>ping</code> | double | Data node response in milliseconds during speed test |

Examples

```
get_data_node()
```

| | |
|-------------------------------|---|
| <code>init_cmip6_index</code> | <i>Build CMIP6 experiment output file index</i> |
|-------------------------------|---|

Description

`init_cmip6_index()` will search the CMIP6 model output file using [esgf_query\(\)](#) , return a `data.table::data.table()` containing the actual NetCDF file url to download, and store it into user data directory for future use.

Usage

```
init_cmip6_index(
  activity = "ScenarioMIP",
  variable = c("tas", "tasmax", "tasmin", "hurs", "hursmax", "hursmin", "pr", "rsds",
    "rlds", "psl", "sfcWind", "clt"),
  frequency = "day",
  experiment = c("ssp126", "ssp245", "ssp370", "ssp585"),
  source = c("AWI-CM-1-1-MR", "BCC-CSM2-MR", "CESM2", "CESM2-WACCM", "EC-Earth3",
    "EC-Earth3-Veg", "GFDL-ESM4", "INM-CM4-8", "INM-CM5-0", "MPI-ESM1-2-HR",
    "MRI-ESM2-0"),
  variant = "r1i1p1f1",
  replica = FALSE,
  latest = TRUE,
  resolution = c("100 km", "50 km"),
  limit = 10000L,
  data_node = NULL,
  years = NULL,
  save = FALSE
)
```

Arguments

| | |
|----------|---|
| activity | A character vector indicating activity identifiers. Default: "ScenarioMIP". Possible values: |
| | <ul style="list-style-type: none"> • "AerChemMIP": Aerosols and Chemistry Model Intercomparison Project, • "C4MIP": Coupled Climate Carbon Cycle Model Intercomparison Project, • "CDRMIP": Carbon Dioxide Removal Model Intercomparison Project, • "CFMIP": Cloud Feedback Model Intercomparison Project, • "CMIP": CMIP DECK: 1pctCO2, abrupt4xCO2, amip, esm-piControl, esm-historical, historical, and piControl experiments, • "CORDEX": Coordinated Regional Climate Downscaling Experiment, • "DAMIP": Detection and Attribution Model Intercomparison Project, • "DCPP": Decadal Climate Prediction Project, • "DynVarMIP": Dynamics and Variability Model Intercomparison Project, • "FAFMIP": Flux-Anomaly-Forced Model Intercomparison Project, • "GMMIP": Global Monsoons Model Intercomparison Project, • "GeoMIP": Geoengineering Model Intercomparison Project, • "HighResMIP": High-Resolution Model Intercomparison Project, • "ISMIP6": Ice Sheet Model Intercomparison Project for CMIP6, • "LS3MIP": Land Surface, Snow and Soil Moisture, • "LUMIP": Land-Use Model Intercomparison Project, • "OMIP": Ocean Model Intercomparison Project, • "PAMIP": Polar Amplification Model Intercomparison Project, • "PMIP": Palaeoclimate Modelling Intercomparison Project, • "RFMIP": Radiative Forcing Model Intercomparison Project, |

| | |
|------------|---|
| | <ul style="list-style-type: none"> • "SIMIP": Sea Ice Model Intercomparison Project, • "ScenarioMIP": Scenario Model Intercomparison Project, • "VIACSAB": Vulnerability, Impacts, Adaptation and Climate Services Advisory Board, • "VolMIP": Volcanic Forcings Model Intercomparison Project |
| variable | A character vector indicating variable identifiers. The 12 most related variables for EPW are set as defaults. If NULL, all possible variables are returned. Default: c("tas", "tasmax", "tasmin", "hurs", "hursmax", "hursmin", "psl", "rss", "rls", "sfcWind", "pr", "prf", "cldt"), where: <ul style="list-style-type: none"> • tas: Near-surface (usually, 2 meter) air temperature, units: K. • tasmax: Maximum near-surface (usually, 2 meter) air temperature, units: K. • tasmin: Minimum near-surface (usually, 2 meter) air temperature, units: K. • hurs: Near-surface relative humidity, units: %. • hursmax: Maximum near-surface relative humidity, units: %. • hursmin: Minimum near-surface relative humidity, units: %. • psl: Sea level pressure, units: Pa. • rsds: Surface downwelling shortwave radiation, units: W m-2. • rlds: Surface downwelling longwave radiation, units: W m-2. • sfcWind: Near-surface (usually, 10 meters) wind speed, units: m s-1. • pr: Precipitation, units: kg m-2 s-1. • cldt: Total cloud area fraction for the whole atmospheric column, as seen from the surface or the top of the atmosphere. Units: %. |
| frequency | A character vector of sampling frequency. If NULL, all possible frequencies are returned. Default: "day". Possible values: <ul style="list-style-type: none"> • "1hr": sampled hourly, • "1hrCM": monthly-mean diurnal cycle resolving each day into 1-hour means, • "1hrPt": sampled hourly, at specified time point within an hour, • "3hr": sampled every 3 hours, • "3hrPt": sampled 3 hourly, at specified time point within the time period, • "6hr": sampled every 6 hours, • "6hrPt": sampled 6 hourly, at specified time point within the time period, • "day": daily mean samples, • "dec": decadal mean samples, • "fx": fixed (time invariant) field, • "mon": monthly mean samples, • "monC": monthly climatology computed from monthly mean samples, • "monPt": sampled monthly, at specified time point within the time period, • "subhrPt": sampled sub-hourly, at specified time point within an hour, • "yr": annual mean samples, • "yrPt": sampled yearly, at specified time point within the time period |
| experiment | A character vector indicating root experiment identifiers. The Tier-1 experiment of activity ScenarioMIP are set as defaults. If NULL, all possible experiment are returned. Default: c("ssp126", "ssp245", "ssp370", "ssp585"). |

| | |
|-------------------|---|
| source | A character vector indicating model identifiers. Defaults are set to 11 sources which give outputs of all 4 experiment of activity ScenarioMIP with daily frequency, i.e. "AWI-CM-1-1-MR", "BCC-CSM2-MR", "CESM2", "CESM2-WACCM", "EC-Earth3", "EC-Earth3-Veg", "GFDL-ESM4", "INM-CM4-8", "INM-CM5-0", "MPI-ESM1-2-HR" and "MRI-ESM2-0". If NULL, all possible sources are returned. |
| variant | A character vector indicating label constructed from 4 indices stored as global attributes in format r<k>i<l>p<m>f<n> described below. Default: "r1i1p1f1". If NULL, all possible variants are returned. <ul style="list-style-type: none"> • r: realization_index (<k>) = realization number (integer >0) • i: initialization_index (<l>) = index for variant of initialization method (integer >0) • p: physics_index (<m>) = index for model physics variant (integer >0) • f: forcing_index (<n>) = index for variant of forcing (integer >0) |
| replica | Whether the record is the "master" copy, or a replica. Use FALSE to return only originals and TRUE to return only replicas. Default: FALSE. |
| latest | Whether the record is the latest available version, or a previous version. Use TRUE to return only the latest version of all records and FALSE to return previous versions. Default: FALSE. |
| resolution | A character vector indicating approximate horizontal resolution. Default: c("50 km", "100 km"). If NULL, all possible resolutions are returned. |
| limit | An integer indicating the maximum of matched records to return. Should be <= 10,000. Default: 10000. |
| data_node | A character vector indicating data nodes to be queried. Default to NULL, which means all possible data nodes. |
| years | An integer vector indicating the target years to be include in the data file. All other years will be excluded. If NULL, no subsetting on years will be performed. Default: NULL. |
| save | If TRUE, the results will be saved into user data directory. Default: FALSE. |

Details

For details on where the file index is stored, see [rappdirs::user_data_dir\(\)](#).

Value

A [data.table::data.table](#) with 22 columns:

| No. | Column | Type | Description |
|-----|----------------|-----------|--|
| 1 | file_id | Character | Model output file universal identifier |
| 2 | dataset_id | Character | Dataset universal identifier |
| 3 | mip_era | Character | Activity's associated CMIP cycle. Will always be "CMIP6" |
| 4 | activity_drs | Character | Activity DRS (Data Reference Syntax) |
| 5 | institution_id | Character | Institution identifier |
| 6 | source_id | Character | Model identifier |
| 7 | experiment_id | Character | Root experiment identifier |

| | | | |
|----|--------------------|-----------|--|
| 8 | member_id | Character | A compound construction from sub_experiment_id and variant_label |
| 9 | table_id | Character | Table identifier |
| 10 | frequency | Character | Sampling frequency |
| 11 | grid_label | Character | Grid identifier |
| 12 | version | Character | Approximate date of model output file |
| 13 | nominal_resolution | Character | Approximate horizontal resolution |
| 14 | variable_id | Character | Variable identifier |
| 15 | variable_long_name | Character | Variable long name |
| 16 | variable_units | Character | Units of variable |
| 17 | datetime_start | POSIXct | Start date and time of simulation |
| 18 | datetime_end | POSIXct | End date and time of simulation |
| 19 | file_size | Character | Model output file size in Bytes |
| 20 | data_node | Character | Data node to download the model output file |
| 21 | dataset_pid | Character | A unique string that helps identify the dataset |
| 22 | tracking_id | Character | A unique string that helps identify the output file |

Note

Argument `limit` will only apply to Dataset query. `init_cmip6_index()` will try to get all model output files which match the dataset id.

Examples

```
## Not run:
init_cmip6_index()

## End(Not run)
```

load_cmip6_index
Load previously stored CMIP6 experiment output file index
Description

Load previously stored CMIP6 experiment output file index

Usage

```
load_cmip6_index(force = FALSE)
```

Arguments

| | |
|--------------------|---|
| <code>force</code> | If TRUE, read the index file. Otherwise, return the cached index if exists. Default: FALSE. |
|--------------------|---|

Value

A `data.table::data.table` with 20 columns. For detail description on column, see [init_cmip6_index\(\)](#).

Examples

```
## Not run:
load_cmip6_index()

## End(Not run)
```

match_coord

Match coordinates of input EPW in the CMIP6 output file database

Description

`match_coord()` takes an EPW and uses its longitude and latitude to match corresponding values that meet specified threshold in NetCDF files.

Usage

```
match_coord(epw, threshold = list(lon = 1, lat = 1), max_num = NULL)
```

Arguments

| | |
|------------------------|---|
| <code>epw</code> | Possible values: |
| | <ul style="list-style-type: none"> • A file path of EPW file • An eplusr::Epw object • A regular expression used to search locations in EnergyPlus Weather Database, e.g. "los angeles.*tmy3". You will be asked to select a matched EPW to download and read. It will be saved into tempdir(). Note that the search is case-insensitive |
| <code>threshold</code> | A list of 2 elements <code>lon</code> and <code>lat</code> specifying the absolute distance threshold used when matching longitude and latitude. Default: <code>list(lon = 1.0, lat = 1.0)</code> |
| <code>max_num</code> | The maximum number to be matched for both longitude and latitude when threshold is matched. Default is <code>NULL</code> , which means no limit |

Details

`match_coord()` uses [future.apply](#) underneath. You can use your preferable [future](#) backend to speed up data extraction in parallel. By default, `match_coord()` uses [future::sequential](#) backend, which runs things in sequential.

Value

An `epw_cmip6_coord` object, which is basically a list of 3 elements:

- `epw`: An [eplusr::Epw](#) object parsed from input `epw` argument
- `meta`: A list containing basic meta data of input EPW, including `city`, `state_province`, `country`, `latitude` and `longitude`.

- coord: A `data.table::data.table()` which is basically CMIP6 index database with an appending new list column coord that contains matched latitudes and longitudes in each NetCDF file. Each element in coord contains 2 elements lat and lon, in which contains the 4 components describing the matched coordinates.
 - index: the indices of matched coordinates
 - value: the actual longitude or latitude in the NetCDF coordinate grids
 - dis: the distance between the coordinate values in NetCDF and input EPW
 - which: The value indices of longitude or latitude in the NetCDF coordinate grids. These values are used to extract the corresponding variable values

Examples

```
## Not run:
# download an EPW from EnergyPlus website
epw <- eplusr::download_weather("los angeles.*TMY3", dir = tempdir(),
  type = "EPW", ask = FALSE)

match_coord(epw, threshold = list(lon = 1.0, lat = 1.0))

## End(Not run)
```

morphing_epw

Morphing EPW weather variables

Description

`morphing_epw()` takes an `epw_cmip6_data` object generated using `extract_data()` and calculates future core EPW weather variables using Morphing Method.

Usage

```
morphing_epw(data, years = NULL, labels = NULL)
```

Arguments

| | |
|---------------------|--|
| <code>data</code> | An <code>epw_cmip6_data</code> object generated using <code>extract_data()</code> |
| <code>years</code> | An integer vector indicating the target years to be considered. If <code>NULL</code> , all years in input data will be considered. Default: <code>NULL</code> . |
| <code>labels</code> | A character or factor vector used for grouping input years. Usually are the outputs of <code>base::cut()</code> . <code>labels</code> should have the same length as <code>years</code> . If given, climate data of <code>years</code> grouped by <code>labels</code> will be averaged. Default: <code>NULL</code> . |

Details

The EPW weather variables that get morphed are listed in details.

Value

An epw_cmip6_morphed object, which is basically a list of 12 elements:

| No. | Element | Type | Morphing Method | Description |
|-----|--------------|--------------------------|-----------------|---|
| 1 | epw | eplusr::Epw | N/A | The original EPW file used for morphing |
| 2 | tdb | data.table::data.table() | Stretch | Data of dry-bulb temperature after morphing |
| 3 | tdew | data.table::data.table() | Derived | Data of dew-point temperature after morphing |
| 4 | rh | data.table::data.table() | Stretch | Data of relative humidity after morphing |
| 5 | p | data.table::data.table() | Stretch | Data of atmospheric pressure after morphing |
| 6 | hor_ir | data.table::data.table() | Stretch | Data of horizontal infrared radiation from the sky |
| 7 | glob_rad | data.table::data.table() | Stretch | Data of global horizontal radiation after morphing |
| 8 | norm_rad | data.table::data.table() | Derived | Data of direct normal radiation after morphing |
| 9 | diff_rad | data.table::data.table() | Stretch | Data of diffuse horizontal radiation after morphing |
| 10 | wind | data.table::data.table() | Stretch | Data of wind speed after morphing |
| 11 | total_cover | data.table::data.table() | Derived | Data of total sky cover after morphing |
| 12 | opaque_cover | data.table::data.table() | Derived | Data of opaque sky cover after morphing |

Each `data.table::data.table()` listed above contains x columns

| No. | Column | Type | Description |
|-----|----------------------|-----------|--|
| 1 | activity_drs | Character | Activity DRS (Data Reference Syntax) |
| 2 | institution_id | Character | Institution identifier |
| 3 | source_id | Character | Model identifier |
| 4 | experiment_id | Character | Root experiment identifier |
| 5 | member_id | Character | A compound construction from sub_experiment_id and variant_label |
| 6 | table_id | Character | Table identifier |
| 7 | lat | Double | The averaged values of input latitude |
| 8 | lon | Double | The averaged values of input longitude |
| 9 | interval | Factor | The year value of data morphed |
| 10 | Variable Name | Double | The morphed data, where Variable Name is the corresponding EPW weather variable name |
| 11 | delta | Double | The shift factor. Will be NA for derived values |
| 12 | alpha | Double | The stretch factor. Will be NA for derived values |

References

Belcher, S., Hacker, J., Powell, D., 2005. Constructing design weather data for future climates. Building Services Engineering Research and Technology 26, 49–61. <https://doi.org/10.1191/0143624405bt112oa>

Description

`set_cmip6_index()` takes a `data.table::data.table()` as input and set it as current index.

Usage

```
set_cmip6_index(index, save = FALSE)
```

Arguments

- index A `data.table::data.table()` containing the same column names and types as the output of `init_cmip6_index()`.
- save If TRUE, besides loaded index, the index file saved to data directory will be also updated. Default: FALSE.

Details

`set_cmip6_index()` is useful when `init_cmip6_index()` may give you too much cases of which only some are of interest.

Value

A `data.table::data.table()`.

summary_database

*Summary CMIP6 model output file status***Description**

`summary_database()` scan the directory specified and returns a `data.table()` containing summary information about all the CMIP6 files available against the output file index loaded using `load_cmip6_index()`.

Usage

```
summary_database(
  dir,
  by = c("activity", "experiment", "variant", "frequency", "variable", "source",
        "resolution"),
  mult = c("skip", "latest"),
  append = FALSE,
  recursive = FALSE,
  update = FALSE,
  warning = TRUE
)
```

Arguments

| | |
|------------------------|---|
| <code>dir</code> | A single string indicating the directory where CMIP6 model output NetCDF files are stored. |
| <code>by</code> | The grouping column to summary the database status. Should be a subset of: <ul style="list-style-type: none"> • "experiment": root experiment identifiers • "source": model identifiers • "variable": variable identifiers • "activity": activity identifiers • "frequency": sampling frequency • "variant": variant label • "resolution": approximate horizontal resolution |
| <code>mult</code> | Actions when multiple files match a same case in the CMIP6 index. If "latest", the file with latest modification time will be used. If "skip", all matched files will be skip and this case will be kept as unmatched. Default: "skip". |
| <code>append</code> | If TRUE, status of CMIP6 files will only be updated if they are not found in previous summary. This is useful if CMIP6 files are stored in different directories. Default: FALSE. |
| <code>recursive</code> | If TRUE, scan recursively into directories. Default: FALSE. |
| <code>update</code> | If TRUE, the output file index will be updated based on the matched NetCDF files in specified directory. If FALSE, only current loaded index will be updated, but the actual index database file saved in <code>get_data_dir()</code> will remain unchanged. Default: FALSE. |
| <code>warning</code> | If TRUE, warning messages will show when multiple files match a same case. Default: TRUE. |

Details

`summary_database()` uses `future.apply` underneath. You can use your preferable `future` backend to speed up data extraction in parallel. By default, `summary_database()` uses `future::sequential` backend, which runs things in sequential.

Value

A `data.table::data.table()` containing corresponding grouping columns plus:

| Column | Type | Description |
|-----------------------------|----------------------|-------------------------------------|
| <code>datetime_start</code> | <code>POSIXct</code> | Start date and time of simulation |
| <code>datetime_end</code> | <code>POSIXct</code> | End date and time of simulation |
| <code>file_num</code> | <code>Integer</code> | Total number of file per group |
| <code>file_size</code> | Units (Mbytes) | Approximate total size of file |
| <code>dl_num</code> | <code>Integer</code> | Total number of file downloaded |
| <code>dl_percent</code> | Units (%) | Total percentage of file downloaded |
| <code>dl_size</code> | Units (Mbytes) | Total size of file downloaded |

Also an attribute `not_matched` is added to the returned `data.table::data.table()` which contains meta data for those CMIP6 output files that are not covered by current CMIP6 output file index.

For the meaning of grouping columns, see `init_cmip6_index()`.

Examples

```
## Not run:  
summary_database()  
  
summary_database(by = "experiment")  
  
## End(Not run)
```

Index

base::cut(), [17](#)
data.table(), [19](#)
data.table::data.table, [6](#), [14](#), [15](#)
data.table::data.table(), [8](#), [11](#), [17–21](#)

eplusr::Epw, [8](#), [10](#), [16](#), [18](#)
epwshift (epwshift-package), [2](#)
epwshift-package, [2](#)
esgf_query, [3](#)
esgf_query(), [11](#)
extract_data, [7](#)
extract_data(), [17](#)

fst::write_fst(), [8](#)
future, [8](#), [16](#), [20](#)
future.apply, [8](#), [16](#), [20](#)
future::sequential, [8](#), [16](#), [20](#)
future_epw, [9](#)

get_data_dir, [10](#)
get_data_dir(), [20](#)
get_data_node, [11](#)

init_cmip6_index, [11](#)
init_cmip6_index(), [15](#), [19](#), [21](#)

load_cmip6_index, [15](#)
load_cmip6_index(), [19](#)

match_coord, [16](#)
match_coord(), [7](#), [8](#)
morphing_epw, [17](#)
morphing_epw(), [9](#)

pingr::ping(), [11](#)

rappdirs::user_data_dir(), [10](#), [14](#)

set_cmip6_index, [18](#)
summary_database, [19](#)