



EnMAP
Technical Report

EnGeoMAP
Tutorial for Application: Basic minerals
and rare earth elements mapping

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15.04.2016



Recommended citation:

Boesche, N.; Mielke, C.; Rogass, C. (2016) EnGeoMAP - Tutorial for Application: Basic minerals and rare earth elements mapping. EnMAP Technical Report, GFZ Data Services.

DOI: <http://doi.org/10.2312.enmap.2016.003>

Imprint

EnMAP Consortium

GFZ Data Services

Telegrafenberg
D-14473 Potsdam

Published in Potsdam, Germany
April 2016

<http://doi.org/10.2312.enmap.2016.003>



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Last manual update

15.04.2016

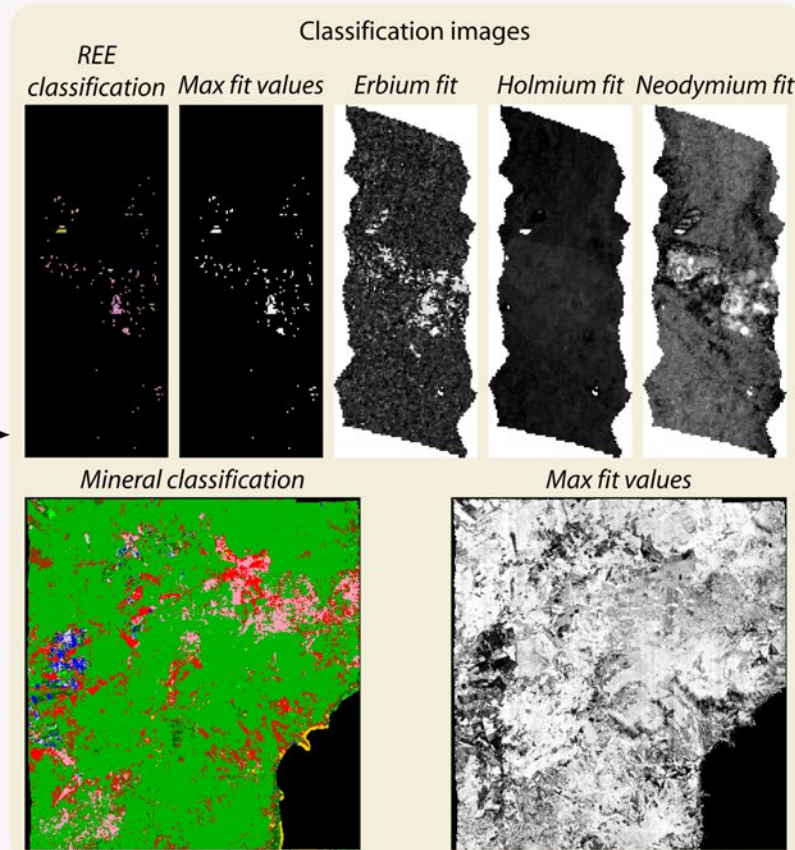
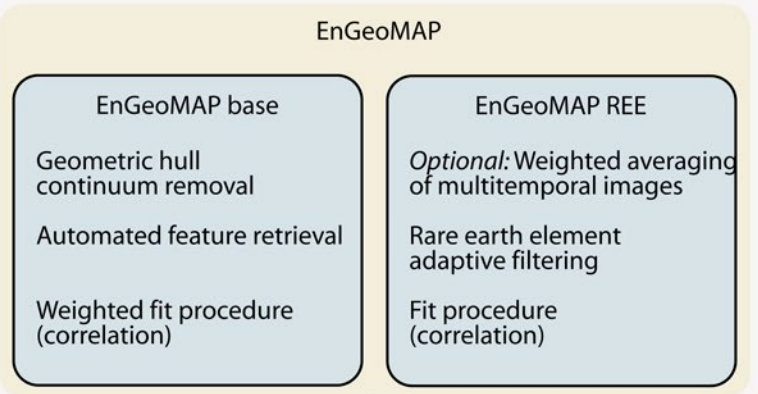
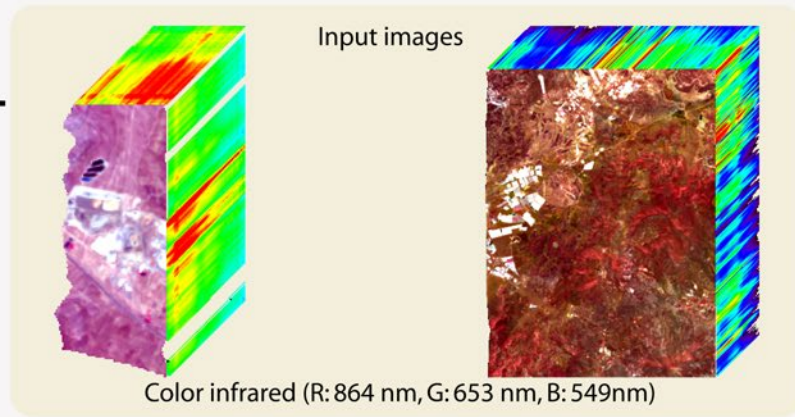
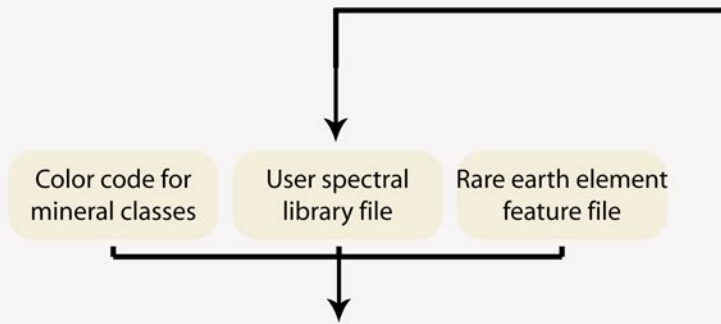
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1 Concept of EnGeoMAP



EnGeoMAP is a program to perform mineral mapping and image classification for geological applications. It contains two sub-programs of particular emphasis on resources determination, mining controlling and environmental protection. These are a basic mineral mapping (EnGeoMAP Base) and a rare earth element mapping (EnGeoMAP REE). In principle, the EnGeoMAP Base program can be used to map Al-OH, Ca-O and Fe-O containing minerals. Rare earth elements can be mapped directly in the surface layer using the EnGeoMAP REE program. Both EnGeoMAP programs utilize a mineral resource type specific spectral library to perform a correlation with the image spectra.

The spectral library for the EnGeoMAP Base program mainly contains a variety of spectra suitable to classify hydrothermal deposits, such as epigenetic and sedimentary hydrothermal deposits (copper, gold, iron, lead, zinc bearing). The spectral library for the EnGeoMAP REE program contains synthetic produced spectra of around 1% of a specific rare earth element in a calcitic matrix. In addition, user specific spectral libraries can be imported into both EnGeoMAP programs. The main difference between both EnGeoMAP programs is the pre-processing of the spectral libraries and satellite input images, as well as the absorption feature definition. EnGeoMAP Base retrieves the feature information automatically by automatic continuum removal (geometric hull) and a signal-to-noise ratio adapted feature definition and retrieval (Mielke et al., 2016). Subsequently, the extracted features are weighted according to their shape and handed over to the correlation step.

The EnGeoMAP REE program consists of a signal-to-noise-ratio weighted averaging of multitemporal acquired images and an adaptive filtering for the rare earth element related absorption bands (Boesche et al., 2015a, Boesche et al., 2015b). In this program the feature retrieval is knowledge based and can be modified by the user.

2 Background

Minerals, or more specifically, ore minerals contain economically valuable elements (mainly metals), which are essential to modern industry and therefore to the development of society. The constantly changing demand of ores and the criticality to the producing industry causes perennial re-evaluations of existing occurrences, deposits and mines, and the global detection and validation of new deposits. Among geophysical and geochemical field surveys, hyperspectral remote sensing surveys are increasingly applied to exploration investigations.

The basic principle of hyperspectral remote sensing for mineral detection is the mathematical description and statistical analysis of material characteristic signals in the spectral ranges of the visible and infrared wavelength range (Clark et al., 2003; Hunt, 1977; Mielke et al., 2016; Swayze et al., 2014; van der Meer, 2004; van der Meer et al., 2012). These characteristic signals (absorption bands) are physically based on electronic transitions in the d- or f-orbitals of the elements (e.g. transitions into the valence band for Fe, and crystal field transitions for rare earth elements), or vibrational motions and their overtones in the molecular bonds (e.g. Fe-OH and Mg-OH in amphiboles, Al-OH in clay minerals, CO₃ in calcium carbonates) (Dieke & Crosswhite, 1963; Hunt, 1977; Mielke et al., 2016; Swayze et al., 2014).

Mineral mapping programs (e.g. the USGS Tetracorder and its successor MICA, and the EnGeoMAP) and image analysis software (e.g. EnMAP-Box, and ENVI-Exelis Visual Information Solutions) utilize the mineral characteristic absorption bands to spectroscopically characterize and map surface materials (Clark et al., 2003; Mielke et al., 2016; van der Linden et al., 2015). The characteristic absorption bands of hyperspectral identifiable minerals serve as a proxy for the lithological units of the different deposit types. These proxy minerals include for example, alunite, chlorite, dickite, epidote, jarosite, kaolinite, and sericite. They serve as key indicators representative for epigenetic and sedimentary hydrothermal deposits (copper, gold, iron, lead, zinc bearing) (Mielke et al., 2016; Swayze et al., 2014).

Carbonatite-alkaline igneous related deposits (rare earth element, lithium, tantalum, niobium bearing) can be mapped using the ankerite, calcite, dolomite, epidote, rare earth element absorption bands, and the clay mineral/carbonates ratio (Boesche, 2015; Boesche et al., 2015a; 2015b; D. J. Turner et al, 2014a; 2014b). In this manner proxy minerals can be found for most deposit types and thus their occurrences can be mapped. In addition to mineral mapping for resource detection, mining monitoring, natural erosion and environmental pollution as well as Earth's crust forming processes (e.g. volcanic activities) can be assessed using the aforementioned mineral mapping methodologies.

3 User Guide

Data import

The EnGeoMAP contains two basic approaches: 1) Base mineral mapping (EnGeoMAP Base) and 2) rare earth element mapping (EnGeoMAP REE).

- Both approaches utilize a spectral library as reference data. The EnGeoMAP Base additionally requires a color codes definition csv-file and EnGeoMAP REE additionally requires a feature definition csv-file (including the name; index-number for Light REE and Heavy REE; the feature shoulder wavelength position; sigma; and the Gaussian-filter width).

Reference files are available in the following directory:

```
EnMAP-Box\SourceCode\applications\EnGeoMAP\_resource
```

```
EnGeoMAP - Base: GFZ_LIB_reduced_EnMAP.esl,  
GFZ_LIB_reduced_EnMAP.hdr, GFZ_LIB_lookup.csv
```

```
EnGeoMAP - REE: GFZLIB_REE_EnMAP, GFZLIB_REE_EnMAP.hdr  
GFZLIB_REE_EnMAP_features.csv
```

A test dataset of Simulated EnMAP Satellite Data for Mountain Pass, USA and Rodalquilar, Spain, is available via the GFZ Data Repository:

<http://doi.org/10.5880/enmap.2016.001>

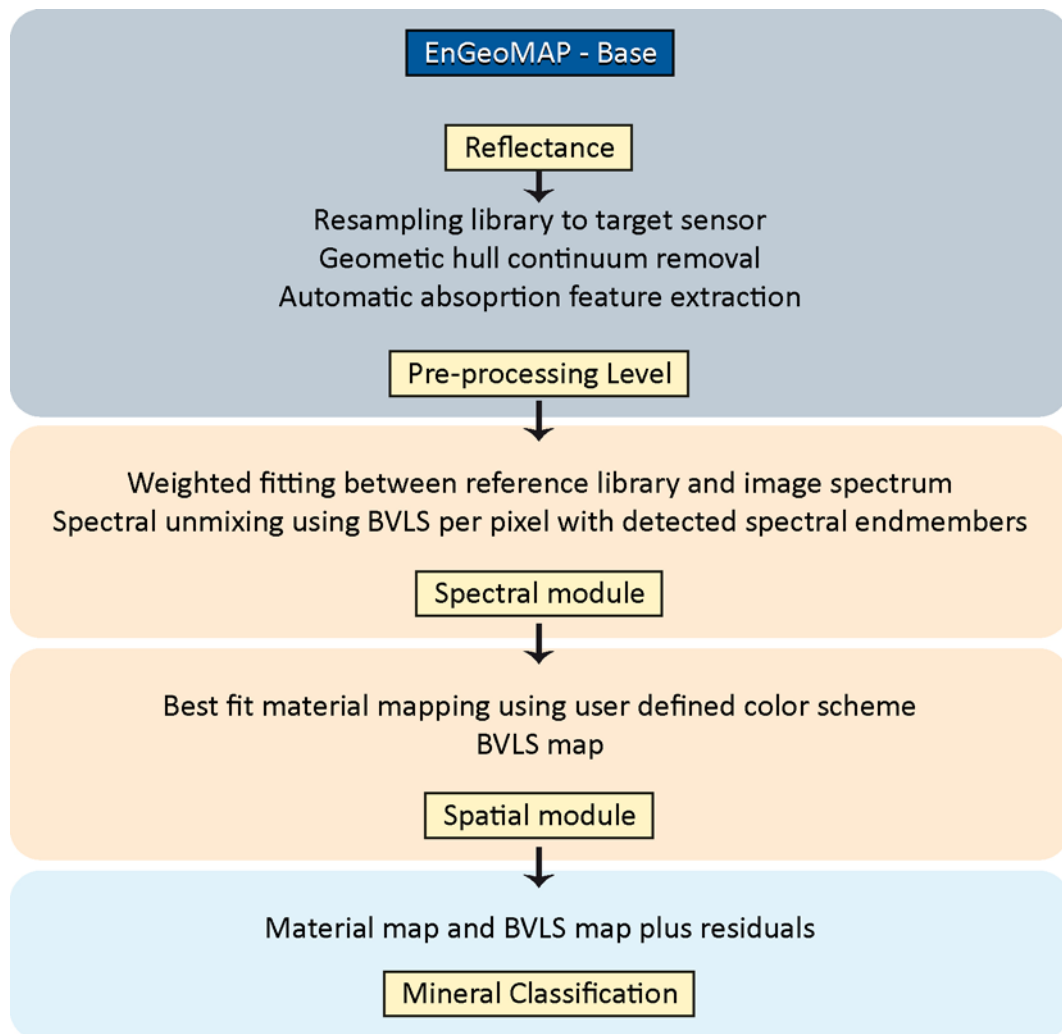
They are freely available to the scientific community under a Creative Commons Attribution-ShareAlike 4.0 International License (Boesche et al., 2016).

When using the data, please cite them as:

Boesche, Nina K.; Mielke, Christian; Segl, Karl; Chabrillat, Sabine; Rogass, Christian; Thomson, David; Lundeen, Sarah; Brell, Maximilian; Guanter, Luis (2016): Simulated EnMAP Satellite Data for Mountain Pass, USA and Rodalquilar, Spain. GFZ Data Services.
<http://doi.org/10.5880/enmap.2016.001>

- The spectral library and the input image files must have the same spectral resolution (example data: EnMAP band centers and band passes). The wavelength start and end value of the spectral library and the images must be the same.
- If multitemporal averaging is used for the EnGeoMAP REE: all input images must be rectified, georeferenced, the subsets identical in size, and must cover the same geographical region. The image data format must be band-sequential (bsq) and the filename extension *.bsq. The image must not contain black background pixel.
- Please do not run ENVI and the EnMAPBOX in one IDL session, as routines may interfere.

EnGeoMAP Base

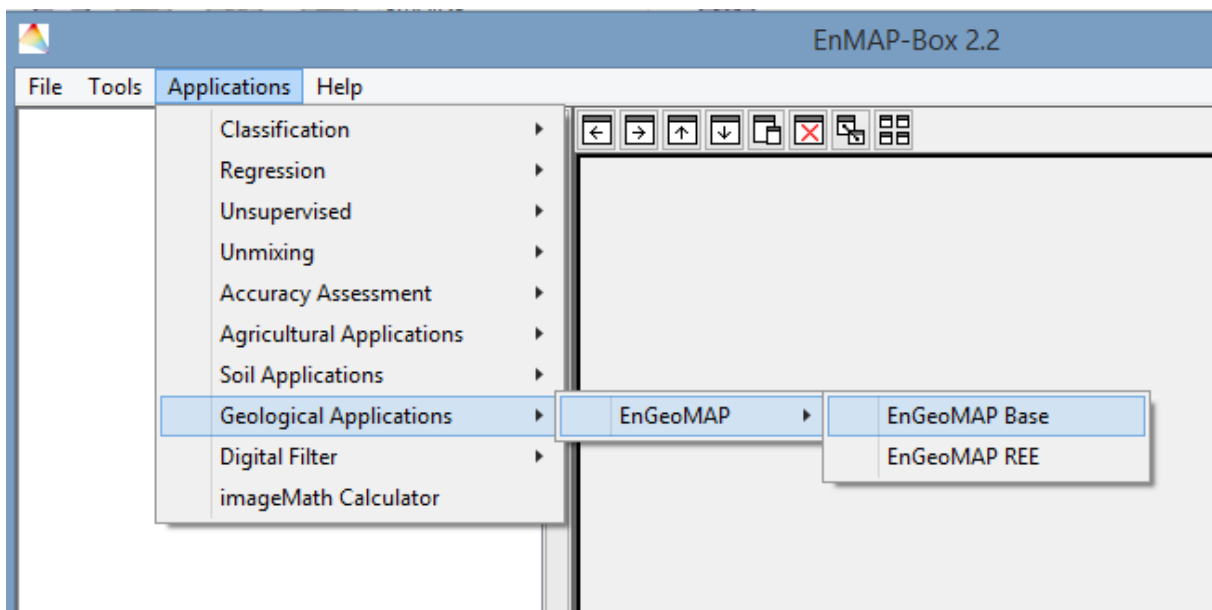


The EnGeoMAP Base program was developed as application for mapping the surface mineralogy from EnMAP. It is based on the geometric hull technique (Mielke et al., 2015; 2016), which allows the extraction of absorption features without any a priori knowledge on absorption feature positions, shape, or the overall shape of the reflectance spectrum itself. The feature definition can either be calculated using the original EnMAP spectral resolution (1 nm precision=No => Faster, Default Option) or a generic 1 nm interpolated spectrum (1 nm precision=Yes => Slower Option).

Users are additionally encouraged to exclude any characteristic absorption features by setting the left and right shoulder positions of the atmospheric water vapor absorption features in the "Exclusion of Atmospheric absorption features" dialogue. Otherwise absorption features that fall within these ranges will also be used in the weighted fitting process! The feature definitions are then used to compute the feature area, feature depth and feature width of the reference features as weights for both, the unknown image spectrum and the library reference spectra. Hence, EnGeoMAP Base uses the $\frac{\text{feature depth} \cdot \text{feature area}}{\text{feature width}}$ in the weighted fit procedure.

Next, the modified spectral angle mapper (MSAM) (Oshigami et al., 2013; Rogass et al., 2013) is used for calculating the fit in case of EnGeoMAP. The resultant data are the weighted fit values of the image pixel spectrum to the reference library entries. A best weighted fit map is produced with the GFZ spectral reference library (Papenfuß, 2015) using the color codes defined in the GFZ_LIB_lookup.csv, which were derived from the color codes of the USGS MICA application (Kokaly, 2012). Additionally, users may conduct a bounded value least squares unmixing on a per pixel basis using only those material entries of the reference library that were detected in the previous best weighted fit step. This might be useful if small material gradients have to be detected.

- Select **Applications > Geological Applications > EnGeoMAP Base**.



- Choose the **Input Image** and the reference **Library File**. Both files need to be presented in the same spectral resolution.
- Optional: A binary **Mask File** can be supplied e.g. to mask out water bodies or vegetation.
- Choose the **Output File**.
- Choose **Flags**:
 - Choose if the Geometric Hull Continuum Removal should be carried out on a generic 1 nm resolution (Yes => Slower) or at the sensors native spectral resolution (No => Faster). Comment: Sensors with very heterogeneous band passes (e.g. HyMAP) may deliver unsatisfactory results if the geometric hull is not calculated at the 1 nm generic resolution.
 - BVLS unmixing (bounded value constrained linear least-squares minimization) at pixel level may be calculated using only those endmembers that have been detected in the weighted fitting process (Default is off). Comment: The material with the highest bvls score will

likely be the material with the largest absorption bands (e.g. iron bearing minerals or vegetation).

- Choose if the GFZ Library is used (default Yes) then a color coded classification result can be computed Based on the color codes in the accompanying lookup table. Comment: The GFZ_LIB_lookup.csv must be saved in the output folder.
- The **Exclusion of Atmospheric Absorption Features** is recommended by supplying the left and right shoulders of the atmospheric water vapor absorption features. Otherwise features found in these spectral regions by the geometric hull will also be used in the weighted fitting process. If necessary, please also remove all residual atmospheric features that fall in the wavelength range before 1250 nm by using sphere interpolation.
- For spiky image spectra it is recommended to use a smoothing filter prior to EnGEOMAP – Base.
- Choose a weighted **Fitting Threshold** (0-1) in order to exclude low correlation values.

GFZ EnGeoMAP - Base v2.008

Input/output

Data file

Library file

Mask file

Output folder C:\Users\username\AppData\Local\Temp\

Flags

1 nm precision Yes No

Additional Bounded Values Least Squares unmixing? Yes No

GFZ spectral library used? GFZ_LIB_lookup.csv must exist! Yes No

Exclusion of Atmospheric absorption features

Example for left shoulders box: 1300;1750 or -1 to disable

Example for right shoulders box: 1450;2010 or -1 to disable

Left shoulders of excluded atmospheric features nm

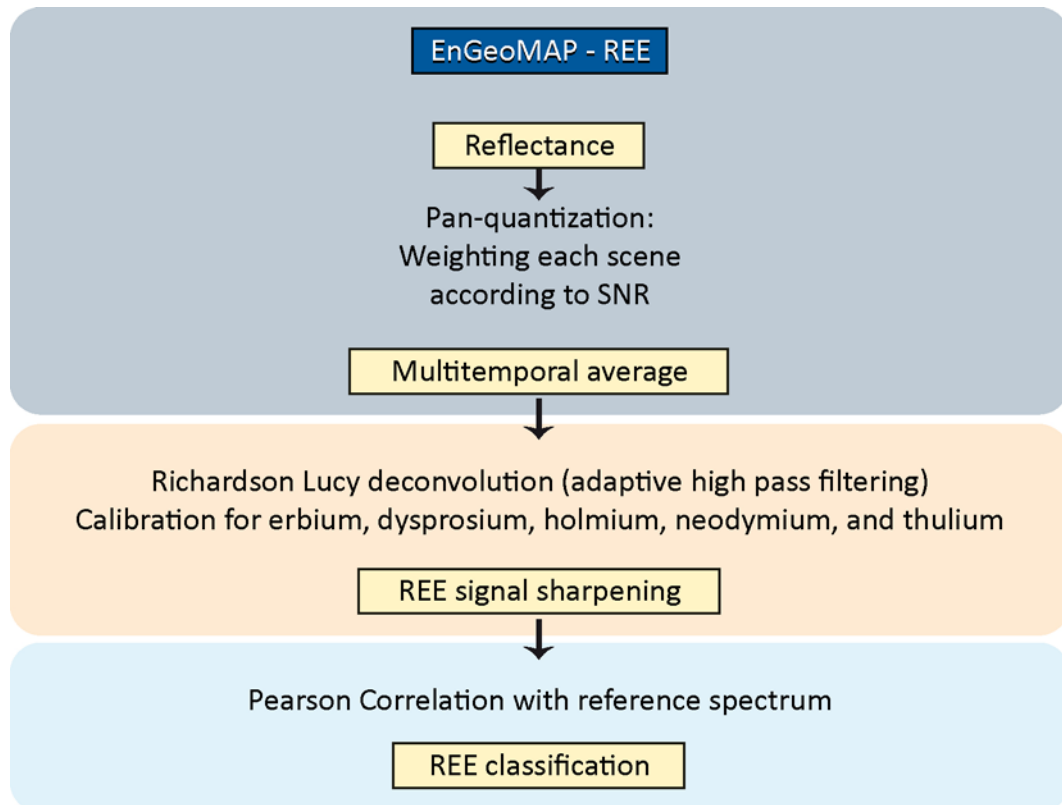
Right shoulders of excluded atmospheric features nm

Fitting threshold

Default 0.0

Accept Cancel

EnGeoMAP REE



Flowchart of EnGEO MAP – REE, after (Nina K. Boesche, 2015)

The EnGeoMAP REE was developed as an additional program to the EnGeoMAP. It separates rare earth element related signatures from noise and absorption features of other minerals. Its advantage is that it finds small absorption features even if they do not dominantly occur as isolated features or are superimposed by noise (Boesche et al., 2015a, Boesche et al., 2015b; Rowan et al., 1986).

The first step of EnGeoMAP REE is a reduction of noise by using a weighted averaging of multi-temporal acquired images. The averaging factor is based on the image signal-to-noise-ratio of homogenous and dark pixels of the input image. These pixels are determined using a “pan-quantization” of the input image (Boesche, 2015). To do that, a spectral subset was extracted that contains only the spectral range covered by the rare earth element that is investigated. The pan image is estimated as a sum over all bands within this subset. Next, the pan image is quantized, which give this step its name. The two first quantiles will then be reduced in size and handed over to determine the signal-to-noise-ratio for low albedos, which serves as the weighting factor for the multitemporal averaging.

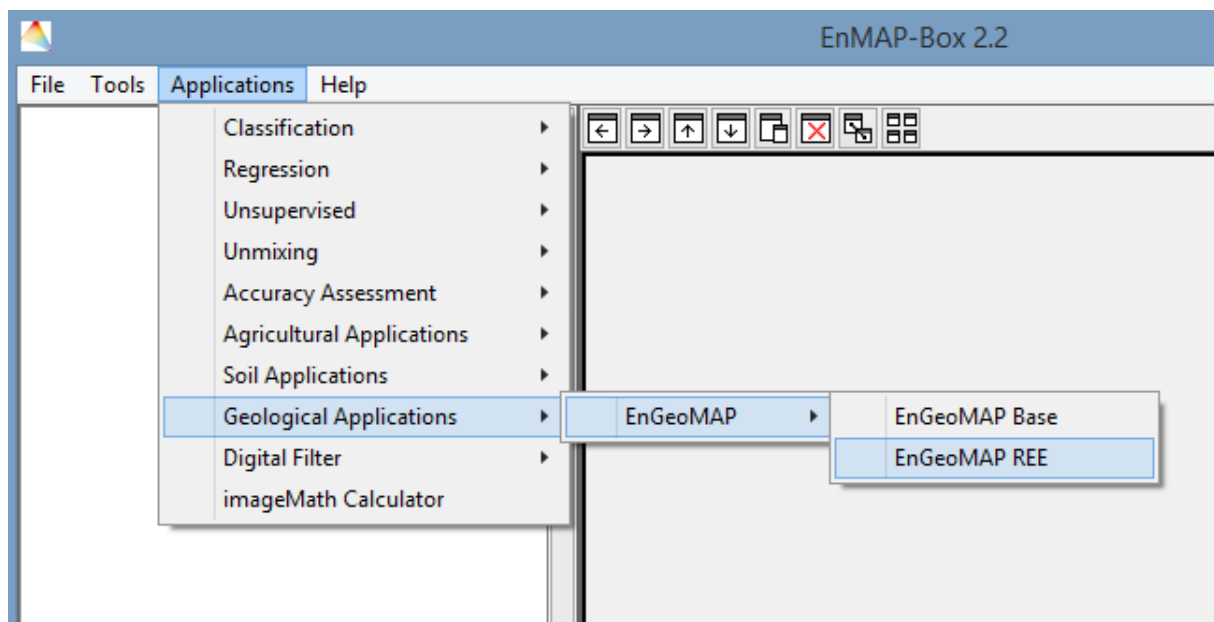
The second step of EnGeoMAP REE is an adaptive filtering of every input image spectrum for rare earth related signals (Boesche, 2015, Boesche et al., 2015a, Boesche et al., 2015b). Filtering is performed using the Richardson-Lucy Deconvolution. For that a filter kernel is defined to perform a two-step deconvolution on the spectrum, see equation (1) and (2). The filter has a Gaussian shape, because it represents the normal distributed probability of orbital energy levels that are involved in the absorption processes (electronic field transitions)

(Sunshine & Pieters, 1993). The Gaussian shape is defined by a sigma value that is a function of the width of the studied rare earth element absorption band. To avoid aliasing, sigma is defined to be at least twice the frequency equivalent to the investigated REE related absorption band.

$$H = H(x, y, \lambda) = \frac{\mathcal{F}\{S^0\}}{\mathcal{F}\{S^i * f\}} f \quad (1)$$

$$S^{i+1} = S^i \circ H(x, y, \lambda) \quad (2)$$

The adaptive filtering is applied to the unknown image spectrum and the library spectrum. In the third and final step, the filtered image spectra are correlated to the filtered spectral library spectra and the fit values saved in the REE_fit_values.bsq file. The REE_max_fit.bsq file contains for every pixel the maximum fit value of all rare earth element fit values. The REE_class.bsq file is a classification file, which contains the assorted rare earth element class for every pixel. The color coding is random and can be modified by the user.



- Select **Applications > Geological Applications > EnGeoMAP REE**.
- Choose the **Input Image**, the reference **Library File** and the **Feature File**. The input image and the reference library file need to be presented in the same spectral resolution. Comment: The output files are saved in the input folder directly.
- Choose **Flag**:
 - Choose if the Rare Earth Element Detection should be carried out on a generic 1 nm resolution (Yes => Slower) or at the sensors native spectral

resolution (No => Faster). Comment: When using the default features that are defined in the GFZLIB_REE_EnMAP_features.csv, sensors with a spectral resolution larger than 4 nm would cause an undersampling in the correlation. In this case EnGeoMAP REE should be carried out on a generic 1 nm resolution.

- The **Exclusion of Atmospheric Absorption Features** is recommended by supplying the left and right shoulders of the atmospheric water vapor absorption features. Otherwise features found in these spectral regions by the geometric hull will also be used in the weighted fitting process.
- Choose a **Fitting Threshold** (0-1) in order to exclude low correlation values. Comment: The default value is 0, therefore all pixel that are positively correlated with the spectral library are flagged to contain a rare earth signal.

GFZ EnGeoMAP - REE v1.007

Input/output

Feature file

Library file

Input/Output folder C:\Users\username\AppData\Local\Temp\

Flags

1 nm precision Yes No

Exclusion of Atmospheric absorption features

Example for left shoulders box: 1300;1750 or -1 to disable
Example for right shoulders box: 1450;2010 or -1 to disable

Left shoulders of excluded atmospheric features nm

Right shoulders of excluded atmospheric features nm

Fitting threshold

Default 0.0

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