Optimality Bounds for a Variational Relaxation of the Image Partitioning Problem

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Abstract We consider a variational convex relaxation of a class of optimal partitioning and multiclass labeling problems, which has recently proven quite successful and can be seen as a continuous analogue of Linear Programming (LP) relaxation methods for finite-dimensional problems. While for the latter case several optimality bounds are known, to our knowledge no such bounds exist in the continuous setting. We provide such a bound by analyzing a probabilistic rounding method, showing that it is possible to obtain an integral solution of the relaxed problem with an a priori upper bound on the objective, ensuring the quality of the result from the viewpoint of optimization. The approach has a natural interpretation as an approximate, multiclass variant of the celebrated coarea formula.

Keywords Convex Relaxation · Multiclass Labeling · Approximation Bound · Combinatorial Optimization · Total Variation · Linear Programming Relaxation

1 Introduction and Background

1.1 Convex Relaxations of Partitioning Problems

In this work, we will be concerned with a class of variational problems used in image processing and analysis for formulating multiclass image partitioning problems, which are of the form

$$\inf_{u \in C} f(u) := \int_{\Omega} (u(x), s(x)) dx + \int_{\Omega} d\Psi(Du),$$

\[ C := \text{BV}(\Omega, \mathcal{E}) \]

with \(\mathcal{E} := \{ e^1, \ldots, e^l \} \).

The labeling function \( u : \Omega \to \mathbb{R}^l \) assigns to each point in the image domain \( \Omega := (0,1)^d \) a label \( i \in I := \{ 1, \ldots, l \} \), which is represented by one of the \( l \)-dimensional unit vectors \( e^1, \ldots, e^l \). Since it is piecewise constant and therefore cannot be assumed to be differentiable, the problem is formulated as a free discontinuity problem in the space \( \text{BV}(\Omega, \mathcal{E}) \) of functions of bounded variation; we refer to [2] for a general overview.

The objective function \( f \) consists of a data term and a regularizer. The data term is given in terms of the \( L^1 \) function \( s(x) = (s_1(x), \ldots, s_l(x)) \in \mathbb{R}^l \), and assigns to the choice \( u(x) = e^i \) the “penalty” \( s_i(x) \), in the sense that

$$\int_{\Omega} (u(x), s(x)) dx = \sum_{i=1}^l \int_{\Omega_i} s_i(x) dx,$$

where \( \Omega_i := u^{-1}(\{e^i\}) = \{ x \in \Omega | u(x) = e^i \} \) is the class region for label \( i \), i.e., the set of points that are assigned the \( i \)-th label. The data term generally depends on the input data – such as color values of a recorded image, depth measurements, or other features – and promotes a good fit of the minimizer to the input data. While it...
is purely local, there are no further restrictions such as continuity, convexity etc., therefore it covers many interesting applications such as segmentation, multi-view reconstruction, stitching, and inpainting [23].

1.2 Convex Regularizers

The regularizer is defined by the positively homogeneous, continuous and convex function \(\Psi: \mathbb{R}^{d \times l} \to \mathbb{R}_{\geq 0}\) acting on the distributional derivative \(Du\) of \(u\), and incorporates additional prior knowledge about the “typical” appearance of the desired output. For piecewise constant \(u\), it can be seen that the definition in (1) amounts to a weighted penalization of the discontinuities of \(u\):

\[
\int_{\Omega} d\Psi(Du) = \int_{J_u} \Psi(\nu_u(x)(u^+(x) - u^-(x))^\top) d\mathcal{H}^{d-1}(x),
\]

where \(J_u\) is the jump set of \(u\), i.e., the set of points where \(u\) has well-defined right-hand and left-hand limits \(u^+\) and \(u^-\) and (in an infinitesimal sense) jumps between the values \(u^+(x), u^-(x) \in \mathbb{R}^d\) across a hyperplane with normal \(\nu_u(x) \in \mathbb{R}^d, \|\nu_u(x)\|_2 = 1\) (see [2] for the precise definitions).

A particular case is to set \(\Psi = (1/\sqrt{2})\|\cdot\|_2\), i.e., the scaled Frobenius norm. In this case \(J(u)\) is just the (scaled) total variation of \(u\) and, since \(u^+(x)\) and \(u^-(x)\) assume values in \(E\) and cannot be equal on the jump set \(J_u\), it holds that

\[
J(u) = \frac{1}{\sqrt{2}} \int_{J_u} \|u^+(x) - u^-(x)\|_2 d\mathcal{H}^{d-1}(x),
\]

\[
= \mathcal{H}^{d-1}(J_u)\tag{7}
\]

Therefore, for \(\Psi = (1/\sqrt{2})\|\cdot\|_2\) the regularizer just amounts to penalizing the total length of the interfaces between class regions as measured by the \((d-1)\)-dimensional Hausdorff measure \(\mathcal{H}^{d-1}\), which is known as uniform metric or Potts regularizer.

A general regularizer was proposed in [18], based on [5]. Given a metric (distance) \(d: \{1, \ldots, l\}^2 \to \mathbb{R}_{\geq 0}\) (not to be confused with the ambient space dimension), define

\[
\Psi_d(z = (z^1, \ldots, z^l)) := \sup_{\nu \in \mathcal{D}_{\text{loc}}} \langle z, v \rangle,
\]

\[
\mathcal{D}_{\text{loc}}^l := \{ (v^1, \ldots, v^l) \in \mathbb{R}^{d \times l} | \ldots \}
\]

\[
\|v^i - v^j\|_2 \leq d(i, j) \forall i, j \in \{1, \ldots, l\}, \ldots
\]

\[
\sum_{k=1}^{l} v^k = 0
\]

It was then shown that

\[
\Psi_d(\nu(e^i - e^j)^\top) = d(i, j),\tag{11}
\]

therefore in view of (7) the corresponding regularizer is non-uniform: the boundary between the class regions \(\Omega_i\) and \(\Omega_j\) is penalized by its length, multiplied by the weight \(d(i, j)\) depending on the labels of both regions.

However, even for the comparatively simple regularizer (4), the model (1) is a (spatially continuous) combinatorial problem due to the integral nature of the constraint set \(C\), therefore optimization is nontrivial. In the context of multiclass image partitioning, a first approach can be found in [20], where the problem was posed in a level set-formulation in terms of a labeling function \(\phi: \Omega \to \{1, \ldots, l\}\), which is subsequently relaxed to \(\mathbb{R}\). Then \(\phi\) is replaced by polynomials in \(\phi\), which coincide with the indicator functions \(u_i\) for the case where \(\phi\) assumes integral values. However, the numerical approach involves several nonlinearities and requires to solve a sequence of nontrivial subproblems.

In contrast, representation (4) directly suggests a more straightforward relaxation to a convex problem: replace \(E\) by its convex hull, which is just the unit simplex in \(l\) dimensions,

\[
\Delta_l := \text{conv}\{e^1, \ldots, e^l\} = \{a \in \mathbb{R}^l | a \geq 0, \sum_{i=1}^{l} a_i = 1\},
\]

and solve the relaxed problem

\[
\inf_{u \in C} f(u),
\]

\[
C := \text{BV}(\Omega, \Delta_l)
\]

\[
= \{ u \in \text{BV}(\Omega)^l \mid u(x) \in \Delta_l \text{ for a.e. } x \in \Omega \}.\tag{15}
\]

Sparked by a series of papers [20,5,17], recently there has been much interest in problems of this form, since they – although generally nonsmooth – are convex and therefore can be solved to global optimality, e.g., using primal-dual techniques. The approach has proven useful for a wide range of applications [13,11,10,28].

1.3 Finite-Dimensional vs. Continuous Approaches

Many of these applications have been tackled before in a finite-dimensional setting, where they can be formulated as combinatorial problems on a grid graph, and solved using combinatorial optimization methods such as \(\alpha\)-expansion and related integer linear programming (ILP) methods [14,15]. These methods have been shown to yield an integral labeling \(u' \in C\) with the a priori bound

\[
f(u') \leq 2 \max_{\alpha \neq j} d(i, j) \min_{\alpha \neq j} d(i, j) f(u'z),
\]

where \(u'z\) is the (unknown) solution of the integral problem (1). They therefore permit to compute a suboptimal solution to the – originally NP-hard – combinatorial problem with an upper bound on the objective.
No such bound is yet available for methods based on the spatially continuous problem \([\Psi]\).

Despite these strong theoretical and practical results available for the finite-dimensional combinatorial energies, the function-based, spatially continuous formulation \([\Psi]\) has several unique advantages:

- The energy \([\Psi]\) is truly isotropic, in the sense that for a proper choice of \(\Psi\) it is invariant under rotation of the coordinate system. Pursuing finite-dimensional “discretize-first” approaches generally introduces artifacts due to the inherent anisotropy, which can only be avoided by increasing the neighborhood size, thereby reducing sparsity and severely slowing down the graph cut-based methods.

In contrast, properly discretizing the relaxed problem \([\Psi]\) and solving it as a convex problem with subsequent thresholding yields much better results without compromising sparsity (Fig. 1 and 2 \([13]\)) . This can be attributed to the fact that solving the discretized problem as a combinatorial problem in effect discards much of the information about the problem structure that is contained in the nonlinear terms of the discretized objective.

- Present combinatorial optimization methods \([4,15]\) are inherently sequential and difficult to parallelize. On the other hand, parallelizing primal-dual methods for solving the relaxed problem \([13]\) is straightforward, and GPU implementations have been shown to outperform state-of-the-art graph cut methods \([29]\).

- Analyzing the problem in a fully functional-analytic setting gives valuable insight into the problem structure, and is of theoretical interest in itself.

### 1.4 Optimality Bounds

However, one possible drawback of the spatially continuous approach is that the solution of the relaxed problem \([\Psi]\) may assume fractional values, i.e., values in \(\Delta_1 \setminus \mathcal{E}\). Therefore, in applications that require a true partition of \(\Omega\), some rounding process is needed in order to generate an integral labeling \(\bar{u}^*\). This may increase the objective, and lead to a suboptimal solution of the original problem \([\Psi]\).

The regularizer \(\Psi_d\) as defined in \([9]\) enjoys the property that it majorizes all other regularizers that can be written in integral form and satisfy \([11]\). Therefore it is in a sense “optimal”, since it introduces as few fractional solutions as possible. In practice, this forces solutions of the relaxed problem to assume integral values in most points, and rounding is in practice only required in small regions.

However, the rounding step may still increase the objective and generate suboptimal integral solutions. Therefore the question arises whether this approach allows to recover “good” integral solutions of the original problem \([\Psi]\).

In the following, we are concerned with the question whether it is possible to obtain, using the convex relaxation \([13]\), integral solutions with an upper bound on the objective. Specifically, we focus on inequalities of the form

\[
f(\bar{u}^*) \leq (1 + \varepsilon)f(u^*_\varepsilon)
\]

(17)

for some constant \(\varepsilon > 0\), which provide an upper bound on the objective of the rounded integral solution \(\bar{u}^*\) with respect to the objective of the (unknown) optimal integral solution \(u^*_\varepsilon\) of \([\Psi]\). Note that generally it is not possible to show that \([17]\) holds for any \(\varepsilon > 0\). The reverse inequality

\[
f(u^*_\varepsilon) \leq f(\bar{u}^*)
\]

(18)

always holds since \(\bar{u}^* \in \mathcal{C}_\varepsilon\) and \(u^*_\varepsilon\) is an optimal integral solution. The specific form \([17]\) can be attributed to the
alternative interpretation
\[ \frac{f(\tilde{u}^*) - f(u_\varepsilon^*)}{f(u_\varepsilon^*)} \leq \varepsilon, \quad (19) \]

which provides a bound for the relative gap to the optimal objective of the combinatorial problem. Such \( \varepsilon \) can be obtained \textit{a posteriori} by actually computing (or approximating) \( \tilde{u}^* \) and a dual feasible point: Assume that a feasible primal-dual pair \((u, v) \in C \times D\) is known, where \( u \) approximates \( u^* \), and assume that some integral feasible \( \tilde{u} \in C \) has been obtained from \( u \) by a rounding process. Then the pair \((\tilde{u}, v)\) is feasible as well since \( C \subseteq C \), and we obtain an \textit{a posteriori} optimality bound of the form \([16]\) with respect to the optimal integral solution \( u_\varepsilon^* \):
\[ \frac{f(\tilde{u}) - f_\mathcal{D}(u_\varepsilon^*)}{f_\mathcal{D}(u_\varepsilon^*)} \leq \frac{f(\tilde{u}) - f_\mathcal{D}(v)}{f_\mathcal{D}(v)} \leq \frac{f(\tilde{u}) - f_\mathcal{D}(v)}{f_\mathcal{D}(v)} =: \varepsilon'. \quad (20) \]

However, this requires that the the primal and dual objectives \( f \) and \( f_\mathcal{D} \) can be accurately evaluated, and requires to compute a minimizer of the problem for the specific input data, which is generally difficult, especially in the spatially continuous formulation.

In contrast, true \textit{a priori} bounds do not require knowledge of a solution and apply uniformly to all problems of a class, irrespective of the particular input. When considering rounding methods, one generally has to discriminate between

- deterministic vs. probabilistic methods, and
- spatially discrete (finite-dimensional) vs. spatially continuous methods.

Most known \textit{a priori} approximation results only hold in the finite-dimensional setting, and are usually proven using graph-based pairwise formulations. In contrast, we assume an “optimize first” perspective due to the reasons outlined in the introduction. Unfortunately, the proofs for the finite-dimensional results often rely on pointwise arguments that cannot directly be transferred to the continuous setting. Deriving similar results for continuous problems therefore requires considerable additional work.

1.5 Contribution and Main Results

In this work we prove that using the regularizer \([16]\), the \textit{a priori} bound \([14]\) can be carried over to the spatially continuous setting. Preliminary versions of these results with excerpts of the proofs have been announced as conference proceedings \([13]\). We extend these results to provide the exact bound \([16]\), and supply the full proofs.

As the main result, we show that it is possible to construct a rounding method parametrized by a parameter \( \gamma \in \Gamma \), where \( \Gamma \) is an appropriate parameter space:
\[ R : C \times \Gamma \rightarrow C \varepsilon, \]
\[ u \in C \rightarrow \tilde{u}_\gamma := R_\gamma(u) \in C \varepsilon, \]
\[ \text{such that for a suitable probability distribution on } \Gamma, \text{ the following theorem holds for the expectation } \mathbb{E} f(\tilde{u}) := \mathbb{E}_\gamma f(\tilde{u}_\gamma): \]

**Theorem 1** Let \( u \in C, s \in L^1(\Omega)^I, \ s \geq 0, \) and let \( \Psi : \mathbb{R}^{d \times l} \rightarrow \mathbb{R}^{\geq 0} \) be positively homogeneous, convex and continuous. Assume there exists a lower bound \( \lambda_1 > 0 \) such that, for \( z = (z^1, \ldots, z^l) \),
\[ \Psi(z) \geq \lambda_1 \frac{1}{2} \sum_{i=1}^l \|z^i\|_2 \quad \forall z \in \mathbb{R}^{d \times l}, \sum_{i=1}^l z^i = 0. \quad (23) \]
Moreover, assume there exists an upper bound \( \lambda_u < \infty \) such that, for every \( \nu \in \mathbb{R}^l \) satisfying \( \|\nu\|_2 = 1 \),
\[ \Psi(\nu(e^i - e^j)^\top) \leq \lambda_u \quad \forall i, j \in \{1, \ldots, l\}. \quad (24) \]
Then Alg. 1 (see below) generates an integral labeling \( \tilde{u} \in C \varepsilon \) almost surely, and the input
\[ \mathbb{E} f(\tilde{u}) \leq 2 \lambda_u f(u). \quad (25) \]

Note that \( \lambda_u \geq \lambda_1 \) always holds if both are defined, since \([23]\) implies, for \( \nu \) with \( \|\nu\|_2 = 1 \),
\[ \lambda_u \geq \Psi(\nu(e^i - e^j)^\top) \geq \frac{\lambda_1}{2} (\|\nu\|_2 + \|\nu\|_2) = \lambda_1. \quad (26) \]

The proof of Thm. 1 (Sect. 4) is based on the work of Kleinberg and Tardos \([12]\), which is set in an LP relaxation framework. However their results are restricted in that they assume a graph-based representation and extensively rely on the finite dimensionality. In contrast, our results hold in the continuous setting without assuming a particular problem discretization.

Theorem 1 guarantees that – in a probabilistic sense – the rounding process may only increase the energy in a controlled way, with an upper bound depending on \( \Psi \). An immediate consequence is

**Corollary 1** Under the conditions of Thm. 1 if \( u^* \) minimizes \( f \) over \( C \), \( u_\varepsilon^* \) minimizes \( f \) over \( C \varepsilon \), and \( u^* \) denotes the output of Alg. 1 applied to \( u^* \), then
\[ \mathbb{E} f(\tilde{u}^*) \leq 2 \frac{\lambda_u}{\lambda_1} f(u_\varepsilon^*). \quad (27) \]

Therefore the proposed approach allows to recover, from the solution \( u^* \) of the convex relaxed problem \([13]\), an approximate integral solution \( \tilde{u}^* \) of the nonconvex
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2 A Probabilistic View of the Coarea Formula

2.1 The Two-Class Case

As a motivation for the following sections, we first provide a probabilistic interpretation of a tool often used in geometric measure theory, the coarea formula (cf. [2]). Assuming $u' \in \text{BV}(\Omega)$ and $u'(x) \in [0, 1]$ for a.e. $x \in \Omega$, the coarea formula states that the total variation of $u$ can be represented by summing the boundary lengths of its super-levelsets:

$$TV(u') = \int_0^1 TV(1_{\{u' > \alpha\}}) d\alpha.$$  \hfill (30)

Here $1_A$ denotes the characteristic function of a set $A$, i.e., $1_A(x) = 1$ iff $x \in A$ and $1_A(x) = 0$ otherwise. The coarea formula provides a connection between problem (11) and the relaxation (13) in the two-class case, where $\mathcal{E} = \{e^1, e^2\}$, $u \in \mathcal{C}_\mathcal{E}$ and $u_1 = 1 - u_2$: As noted in [10],

$$TV(u) = \|e^1 - e^2\|_2 TV(u_1) = \sqrt{2} TV(u_1),$$  \hfill (31)

therefore the coarea formula (30) can be rewritten as

$$TV(u) = \sqrt{2} \int_0^1 TV(1_{\{u_1 > \alpha\}}) d\alpha$$

$$= \int_0^1 TV(e^1 1_{\{u_1 > \alpha\}} + e^2 1_{\{u_1 \leq \alpha\}}) d\alpha$$

$$= \int_0^1 TV(\tilde{u}_\alpha) d\alpha,$$  \hfill (33)

$$\tilde{u}_\alpha := e^1 1_{\{u_1 > \alpha\}} + e^2 1_{\{u_1 \leq \alpha\}}.$$  \hfill (34)

Consequently, the total variation of $u$ can be expressed as the mean over the total variations of a set of integral labelings $\{\tilde{u}_\alpha \in \mathcal{C}_\mathcal{E} | \alpha \in [0, 1]\}$, obtained by rounding $u$ at different thresholds $\alpha$. We now adopt a probabilistic view of (34). We regard the mapping

$$R : (u, \alpha) \in \mathcal{C} \times [0, 1] \mapsto \tilde{u}_\alpha \in \mathcal{C}_\mathcal{E} \quad \text{(a.e. } \alpha \in [0, 1])$$  \hfill (36)

as a parametrized, deterministic rounding algorithm that depends on $u$ and on an additional parameter $\alpha$. From this we obtain a probabilistic (randomized) rounding algorithm by assuming $\alpha$ to be a uniformly distributed random variable. Under these assumptions the coarea formula (34) can be written as

$$TV(u) = \mathbb{E}_\alpha TV(\tilde{u}_\alpha).$$  \hfill (37)

This has the probabilistic interpretation that applying the probabilistic rounding to (arbitrary, but fixed) $u$ does – in a probabilistic sense, i.e., in the mean – not change the objective. It can be shown that this property extends to the full functional $f$ in (13): In the two-class case, the “coarea-like” property

$$f(u) = \mathbb{E}_\alpha f(\tilde{u}_\alpha)$$  \hfill (38)

holds. Functions with property (38) are also known as levelable functions [8,9] or discrete total variations [10] and have been studied in [20]. A well-known implication is that if $u = u^*$, i.e., $u$ minimizes the relaxed problem (13), then in the two-class case almost every $\tilde{u}^* = \tilde{u}_\alpha^*$ is an integral minimizer of the original problem (11), i.e., the optimality bound (17) holds with $\epsilon = 0$ [7].
2.2 The Multi-Class Case and Generalized Coarea Formulas

Generalizing these observations to more than two labels hinges on a property similar to (38) that holds for vector-valued $u$. In a general setting, the question is whether there exist

- a probability space $(\Gamma, \mu)$, and
- a parametrized rounding method, i.e., for $\mu$-almost every $\gamma \in \Gamma$:

$$ R : \mathcal{C} \times \Gamma \to \mathcal{C}_\gamma, \quad u \in \mathcal{C} \mapsto u_{\gamma} := R_{\gamma}(u) \in \mathcal{C}_\gamma $$(39)

satisfying $R_{\gamma}(u') = u'$ for all $u' \in \mathcal{C}_\gamma$, such that a “multiclass coarea-like property” (or generalized coarea formula)

$$ f(u) = \int_{\Gamma} f(\bar{u}_\gamma) d\mu(\gamma) $$ (41)

holds. In a probabilistic sense this corresponds to

$$ f(u) = \int_{\Gamma} f(\bar{u}_\gamma) d\mu(\gamma) = \mathbb{E}_\gamma f(\bar{u}_\gamma), $$ (42)

For $l = 2$ and $\Psi(x) = \| \cdot \|_2$, (31) shows that (42) holds with $\gamma = \alpha$, $\Gamma = [0, 1]$, $\mu = L^1$, and $R : \mathcal{C} \times \Gamma \to \mathcal{C}_\gamma$ as defined in (30). Unfortunately, property (37) is intrinsically restricted to the two-class case with TV regularizer.

In the multiclass case, the difficulty lies in providing a suitable combination of a probability space $(\Gamma, \mu)$ and a parametrized rounding step $(u, \gamma) \mapsto \bar{u}_\gamma$. Unfortunately, obtaining a relation such as (37) for the full functional $\Omega$ is unlikely, as it would mean that solutions to the (after discretization) NP-hard problem (41) could be obtained by solving the convex relaxation (36) and subsequent rounding, which can be achieved in polynomial time.

Therefore we restrict ourselves to an approximate variant of the generalized coarea formula:

$$(1 + \varepsilon) f(u) \geq \int_{\Gamma} f(\bar{u}_\gamma) d\mu(\gamma) = \mathbb{E}_\gamma f(\bar{u}_\gamma). $$ (43)

While (43) is not sufficient to provide a bound on $f(\bar{u}_\gamma)$ for particular $\gamma$, it permits a probabilistic bound for any minimizer $u^*$ of the relaxed problem (13), eq. (43) implies

$$ \mathbb{E}_\gamma f(\bar{u}_\gamma) \leq (1 + \varepsilon) f(u^*) \leq (1 + \varepsilon) f(u^*), $$ (44)

i.e., the ratio between the objective of the rounded relaxed solution and the optimal integral solution is bounded – in a probabilistic sense – by $(1 + \varepsilon)$.

In the following sections we construct a suitable parametrized rounding method and probability space in order to obtain an approximate generalized coarea formula of the form (43).

Algorithm 1 Continuous Probabilistic Rounding

1: $u^0 \leftarrow u$, $U^0 \leftarrow \Theta$, $\varepsilon^0 \leftarrow (1, \ldots, 1) \in \mathbb{R}^l$.
2: for $k = 1, 2, \ldots$ do
3: Randomly choose $\gamma^k = (\bar{i}^k, \alpha^k) \in \mathcal{I} \times [0, 1]$ uniformly.
4: $U^k \leftarrow U^{k-1} \cap \{x \in \Omega | u^k_{\bar{i}^k-1}(x) > \alpha^k\}$.
5: $u^k \leftarrow e^T 1_{M^k} + u^{k-1} 1_{\Theta \setminus M^k}$.
6: $U^k \leftarrow U^{k-1} \setminus M^k$.
7: $c_j^k \leftarrow \min\{e^{k-1}, \alpha^k\}$, $j = \bar{i}^k$.
8: end for

3 Probabilistic Rounding for Multiclass Image Partitions

3.1 Approach

We consider the probabilistic rounding approach based on (12) as defined in Alg. 1.

The algorithm proceeds in a number of phases. At each iteration, a label and a threshold $\gamma^k := (\bar{i}^k, \alpha^k) \in \mathcal{I}^* := \mathcal{I} \times [0, 1]$ are randomly chosen (step 3), and label $\bar{i}^k$ is assigned to all yet unassigned points $x$ where $u^k_{\bar{i}^k-1}(x) > \alpha^k$ holds (step 5). In contrast to the two-class case considered above, the randomness is provided by a sequence $(\gamma^k)$ of uniformly distributed random variables, i.e., $\mathcal{I}^* = (\mathcal{I})^N$.

After iteration $k$, all points in the set $U^k \subseteq \Theta$ are still unassigned, while all points in $\Theta \setminus U^k$ have been assigned an (integral) label in iteration $k$ or in a previous iteration. Iteration $k+1$ potentially modifies points only in the set $U^k$. The variable $c_j^k$ stores the lowest threshold $\alpha$ chosen for label $j$ up to and including iteration $k$, and is only required for the proofs.

While the algorithm is defined using pointwise operations, it is well-defined in the sense that for fixed $\gamma$, the sequence $(u^k)$, viewed as elements in $L^1$, does not depend on the specific representative of $u$ in its equivalence class in $L^1$. The sequences $(M^k)$ and $(U^k)$ depend on the representative, but are unique up to $L^d$-negligible sets.

In an actual implementation, the algorithm could be terminated as soon as all points in $\Theta$ have been assigned a label, i.e., $U^k = \emptyset$. However, in our framework used for analysis the algorithm never terminates explicitly. Instead, for fixed input $u$ we regard the algorithm as a mapping between sequences of parameters (or instances of random variables) $\gamma = (\gamma^k) \in \Gamma$ and sequences of states $(u^k_\gamma)$, $(U^k_\gamma)$ and $(c^k_\gamma)$. We drop the subscript $\gamma$ if it does not create ambiguities. The elements of the sequence $(\gamma^k)$ are independently uniformly distributed, and by the Kolmogorov extension theorem [21] Thm. 2.1.5] there exists a probability space and a stochas-
tic process on the set of sequences $\gamma$ with compatible marginal distributions.

In order to define the parametrized rounding step $(u, \gamma) \mapsto \bar{u}_\gamma$, we observe that once $U_{k'} = \emptyset$ occurs for some $k' \in \mathbb{N}$, the sequence $(u^k_{\gamma})$ becomes stationary at $u^k_{\gamma}$. In this case the algorithm may be terminated, with output $\bar{u}_\gamma := u^k_{\gamma}$.

**Definition 1** Let $u \in BV(\Omega)^l$ and $f : BV(\Omega)^l \rightarrow \mathbb{R}$. For some $\gamma, k \in \Gamma$, if $U_{k'} = \emptyset$ in Alg. 1 for some $k' \in \mathbb{N}$, we denote $\bar{u}_\gamma := u^k_{\gamma}$. We define

$$f(\bar{u}_\gamma) := \{ f(u^k_{\gamma}) : \exists k' \in \mathbb{N} : U_{k'} = \emptyset \land u^k_{\gamma} \in BV(\Omega)^l, \}$$

We denote by $f(\bar{u})$ the corresponding random variable induced by assuming $\gamma$ to be uniformly distributed on $\Gamma$.

As indicated above, $f(\bar{u}_\gamma)$ is well-defined: if $U_{k'} = \emptyset$ for some $(\gamma, k')$ then $u^k_{\gamma} = u^{k''}_{\gamma}$ for all $k'' \geq k'$. Instead of focusing on local properties of the random sequence $(u^k_{\gamma})$ as in the proofs for the finite-dimensional case, we derive our results directly for the sequence $(f(u^k_{\gamma}))$. In particular, we show that the expectation of $f(\bar{u})$ over all sequences $\gamma$ can be bounded according to

$$\mathbb{E}[f(\bar{u})] = \mathbb{E}_\gamma f(\bar{u}_\gamma) \leq (1 + \varepsilon)f(\bar{u})$$

for some $\varepsilon \geq 0$, cf. (48). Consequently, the rounding process may only increase the average objective in a controlled way.

### 3.2 Termination Properties

Theoretically, the algorithm may produce a sequence $(u^k_{\gamma})$ that does not become stationary, or becomes stationary with a solution that is not an element of $BV(\Omega)^l$. In Thm. 2 below we show that this happens only with zero probability, i.e., almost surely Alg. 1 generates in a finite number of iterations an integral labeling function $\bar{u}_\gamma \in C_{\mathcal{E}}$. The following two propositions are required for the proof.

**Proposition 1** For the sequence $(c^k)$ generated by Algorithm 1,

$$\mathbb{P}(c^k \sim c^k < 1) \geq$$

$$\sum_{p \in \{0,1\}^t} (-1)^{\mathbb{E}(p)} \left( \sum_{j=1}^t 1 \left( \frac{1}{7} \right)^{p_j} \right)^k$$

holds. In particular,

$$\mathbb{P}(c^k \sim c^k < 1) \rightarrow \infty.$$

**Proof** Denote by $n^k_j \in \mathbb{N}_0$ the number of $k' \in \{1, \ldots, k\}$ such that $i_{k'} = j$, i.e., the number of times label $j$ was selected up to and including the $k$-th step. Then

$$\mathbb{P}(n^k_1, \ldots, n^k_k) \sim \text{Multinomial} \left( k; \frac{1}{7}, \ldots, \frac{1}{7} \right),$$

i.e., the probability of a specific instance is

$$\mathbb{P}(n^k_1, \ldots, n^k_k) = \frac{k!}{n^k_1! \cdots n^k_k!} \left( \frac{1}{7} \right)^k \cdot \sum_{n^k_1 = k} \mathbb{P}(c^k \sim c^k < 1 | (n^k_1, \ldots, n^k_k)).$$

Therefore,

$$\mathbb{P}(c^k \sim c^k < 1) \geq \sum_{n^k_1 + \cdots + n^k_k = k} \frac{k!}{n^k_1! \cdots n^k_k!} \left( \frac{1}{7} \right)^k \cdot \mathbb{P}(c^k \sim c^k < 1 | (n^k_1, \ldots, n^k_k)).$$

Since $c^k_0, \ldots, c^k_k \sim p$ is a sufficient condition for $c^k \sim c^k < 1$, we may bound the probability according to

$$\mathbb{P}(c^k \sim c^k < 1) \geq \sum_{n^k_1 + \cdots + n^k_k = k} \frac{k!}{n^k_1! \cdots n^k_k!} \left( \frac{1}{7} \right)^k \cdot \mathbb{P} \left( c^k \sim c^k < 1 \& j \in \mathcal{I}|(n^k_1, \ldots, n^k_k) \right).$$

We now consider the distributions of the components $c^k_j$ of $c^k$ conditioned on the vector $(n^k_1, \ldots, n^k_k)$. Given $n^k_j$, the probability of $\{c^k_j \geq t\}$ is the probability that in each of the $n^k_j$ steps where label $j$ was selected the threshold $\alpha$ was randomly chosen to be at least as large as $t$. For $0 < t < 1$, we conclude

$$\mathbb{P}(c^k_j \sim c^k_j < t | (n^k_1, \ldots, n^k_k)) = \mathbb{P}(c^k < t | n^k_j) \geq 1 - \mathbb{P}(c^k_j \geq t | n^k_j) \geq 1 - (1 - t)^{n^k_j}.$$  

The above formulation also covers the case $n^k_j = 0$ (note that we assumed $0 < t < 1$). For fixed $k$ the distributions of the $c^k_j$ are independent when conditioned on $(n^k_1, \ldots, n^k_k)$. Therefore we obtain from (53) and (54)

$$\mathbb{P}(c^k \sim c^k < 1) \geq \sum_{n^k_1 + \cdots + n^k_k = k} \frac{k!}{n^k_1! \cdots n^k_k!} \left( \frac{1}{7} \right)^k \cdot \prod_{j=1}^t \mathbb{P} \left( c^k_j \sim c^k_j < t | (n^k_1, \ldots, n^k_k) \right),$$

$$\prod_{j=1}^t \mathbb{P} \left( c^k_j \sim c^k_j < t | (n^k_1, \ldots, n^k_k) \right).$$

(57) and (58)
Expanding the product and swapping the summation order, we derive
\[
\begin{align*}
\mathbb{P}(e^T c_k < 1) & \geq \sum_{n_1 + \ldots + n_l = k} \frac{k!}{n_1! \ldots n_l!} \left( \frac{1}{l} \right)^k \\
& \quad \times \prod_{j=1}^l \left( 1 - \frac{1}{l} \right)^{n_j} p_j, \\
& = \sum_{p \in \{0,1\}^l} (-1)^e p \prod_{j=1}^l \left( \frac{1}{l} \right)^{n_j}.
\end{align*}
\]
(60)

Using the multinomial summation formula, we conclude
\[
\mathbb{P}(e^T c_k < 1) \geq \sum_{p \in \{0,1\}^l} (-1)^e p \left( \prod_{j=1}^l \left( 1 - \frac{1}{l} \right)^{n_j} \right)^k,
\]
(62)

which proves (44). At (8) the multinomial summation formula was invoked. Note that in (62) the \(n_j\) do not occur explicitly anymore. To show the second assertion (48), we use the fact that, for any \(p \neq (0, \ldots, 0)\), \(q_p\) can be bounded by \(0 < q_p < 1\). Therefore
\[
\mathbb{P}(e^T c_k < 1) \geq q_0 + \sum_{p \in \{0,1\}^l, p \neq 0} (-1)^e p (q_p)^k
\]
(63)

\[
= 1 + \sum_{p \in \{0,1\}^l, p \neq 0} (-1)^e p (q_p)^k
\]
(64)

\[
k \leq \sum_{p \in \{0,1\}^l, p \neq 0} (-1)^e p (q_p)^k
\]
(65)

which proves (48).

We now show that Alg. 1 generates a sequence in \(BV(\Omega)^l\) almost surely. The perimeter of a set \(A\) is defined as the total variation of its characteristic function \(\text{Per}(A) := TV(1_A)\).

**Proposition 2** For the sequences \((u^k), (U^k)\) generated by Alg. 1, define
\[
A := \bigcap_{k=1}^\infty \{ \gamma \in \Gamma | \text{Per}(U^k_\gamma) < \infty \}.
\]
(66)

Then
\[
\mathbb{P}(A) = 1.
\]
(67)

If \(\text{Per}(U^k) < \infty\) for all \(k\), then \(u^k \in BV(\Omega)^l\) for all \(k\) as well. Moreover,
\[
\mathbb{P}(u^k \in BV(\Omega)^l \wedge \text{Per}(U^k) < \infty \forall k \in \mathbb{N}) = 1,
\]
(68)

*i.e., the algorithm almost surely generates a sequence of BV functions \((u^k)\) and a sequence of sets of finite perimeter \((U^k)\).*

**Proof** We first show that if \(\text{Per}(U^k) < \infty\) for all \(k' \leq k\), then \(u^k \in BV(\Omega)^l\) for all \(k' \leq k\) as well. For \(k = 0\), the assertion holds, since \(u^0 = u \in BV(\Omega)^l\) by assumption. For \(k \geq 1\),
\[
u^k = e^k 1_{M^k} + u^{k-1} 1_{\Omega \setminus M^k}.
\]
(69)

Since \(M^k = U^{k-1} \cap (\Omega \setminus U^k)\), and \(U^{k}, U^{k-1}\) are assumed to have finite perimeter, \(M^k\) also has finite perimeter. Applying [2, Thm. 3.84] together with the boundedness of \(u^{k-1}\) and \(u^{-1} \in BV(\Omega)^l\) by induction then provides \(u^k \in BV(\Omega)^l\).

We now denote
\[
I^k := \{ \gamma \in \Gamma | \text{Per}(U^k_\gamma) = \infty \},
\]
(70)

and the event that the first set with non-finite perimeter is encountered at step \(k \in \mathbb{N}_0\) by
\[
B_k := I^k \cap (I^k \setminus I^{k-1}) \ldots \cap (I^0 \setminus I^0).
\]
(71)

Then
\[
\mathbb{P}(A) = 1 - \mathbb{P}\left( \bigcup_{k=0}^\infty B_k \right).
\]
(72)

As the sets \(B_k\) are pairwise disjoint, and due to the countable additivity of the probability measure, we have
\[
\mathbb{P}(A) = 1 - \sum_{k=0}^\infty \mathbb{P}(B_k).
\]
(73)

Now \(U^0 = \Omega\), therefore \(\text{Per}(U^0) = TV(1_{\Omega}) = 0 < \infty\) and \(\mathbb{P}(B^0) = 0\). For \(k \geq 1\), we have
\[
\mathbb{P}(B^k) \leq \mathbb{P}\left( \text{Per}(U^k) = \infty \wedge \text{Per}(U^{k'}) < \infty \forall k' < k \right)
\]
\[
\leq \mathbb{P}\left( \text{Per}(U^k) = \infty | \text{Per}(U^{k'}) < \infty \forall k' < k \right)
\]
\[
= \mathbb{P}\left( \text{Per}(U^{k-1}) \cap \{ x \in \Omega | u^{k-1}_x (x) \leq \alpha^k \} \right) = \infty | \text{Per}(U^{k'}) < \infty \forall k' < k \right).
\]
(74)

By the argument from the beginning of the proof, we know that \(u^{k-1} \in BV(\Omega)^l\) under the condition on the perimeter \(\text{Per}(U^{k'})\), therefore from [2, Thm. 3.40] we conclude that \(\text{Per}(\{ x \in \Omega | u^{k-1}_x (x) \leq \alpha^k \})\) is finite for \(L^1\)-a.e. \(\alpha^k\) and all \(k^k\). As the sets of finite perimeter are closed under finite intersection, and since the \(\alpha^k\) are drawn from an uniform distribution, this implies that
\[
\mathbb{P}(\text{Per}(U^k) < \infty | \text{Per}(U^{k-1}) < \infty) = 1.
\]
(75)

Together with (74) we arrive at
\[
\mathbb{P}(B^k) = 0.
\]
(76)

Substituting this result into (72) leads to the assertion,
\[
\mathbb{P}(A) = 1.
\]
(77)

Equation (68) follows immediately.
Proof of the Main Theorem

In order to show the bound and Thm. 1, we first need several technical propositions regarding the composition of two BV functions along a set of finite perimeter. We denote by $(E)^i$ and $(E)^0$ the measure-theoretic interior and exterior of a set $E$, see [2].

$$(E)^i := \{ x \in \Omega : \lim_{\rho \to 0} \frac{|B_\rho(x) \cap E|}{|B_\rho(x)|} = t \}, \ t \in [0, 1].$$

Here $B_\rho(x)$ denotes the ball with radius $\rho$ centered in $x$, and $|A| := \mathcal{L}^d(A)$ the Lebesgue content of a set $A \subseteq \mathbb{R}^d$.

**Proposition 3** Let $\Psi$ be positively homogeneous and convex, and satisfy the upper-boundedness condition [24]. Then

$$(\Psi(\nu(z^1 - z^2) \leq \lambda_u, \ \forall z^1, z^2 \in \Delta_t).$$

Moreover, there exists a constant $C < \infty$ such that

$$\Psi(w) \leq C\|w\|_2 \ \forall w \in W,$$

$$W := \{ w = (w^1 \ldots w^d) \in \mathbb{R}^{d \times l} : \sum_{i=1}^l w^i = 0 \}.$$

**Proof** See appendix.

**Proposition 4** Let $E, F \subseteq \Omega^d$ be $\mathcal{L}^d$-measurable sets. Then

$$(E \cap F)^1 = (E)^1 \cap (F)^1.$$  

**Proof** See appendix.

**Proposition 5** Let $u, v \in BV(\Omega, \Delta_t)$ and $E \subseteq \Omega$ such that $\text{Per}(E) < \infty$. Define

$$w := u1_E + v1_{\Omega \setminus E}.$$  

Then $w \in BV(\Omega, \Delta_t)^1$, and

$$ Dw = Dv_{\perp E}_{\perp E}^1 + Dv_{\perp E}_{\perp E} + \nu E \left( u_{E \setminus F} - v_{E \setminus F} \right) \mathcal{H}^{d-1} (F \cap \Omega),$$

where $u_{E \setminus F}$ and $v_{E \setminus F}$ denote the one-sided approximate limits of $u$ and $v$ on the reduced boundary $F \cap E$, and $\nu E$ is the generalized inner normal of $E$ [2]. Moreover, for continuous, convex and positively homogeneous $\Psi$ satisfying the upper-boundedness condition [24] and some Borel set $A \subseteq \Omega$,

$$\int_A \frac{d\Psi(Dw)}{\mathcal{A}(E)^1} \leq \int_{A \cap (E)^1} \frac{d\Psi(Du)}{\mathcal{A}(E)^1} + \int_{A \cap (E)^0} \frac{d\Psi(Dv)}{\mathcal{A}(E)^1} + \lambda_u \text{Per}(E).$$

**Proof** See appendix.

**Proposition 6** Let $u, v \in BV(\Omega, \Delta_t), \ E \subseteq \Omega$ such that $\text{Per}(E) < \infty$, and

$$u_{|E^1} = v_{|E^1} : \mathcal{L}^d-\text{a.e.}$$

Then $(Du)_{\perp E}_{\perp E}^1 = (Dv)_{\perp E}_{\perp E}^1$, and $\Psi(Du)_{|E^1}^1 = \Psi(Dv)_{|E^1}^1$. In particular,

$$\int_{(E)^1} d\Psi(Du) = \int_{(E)^1} d\Psi(Dv).$$

The result also holds when $(E)^1$ is replaced by $(E)^0$. Moreover, the condition [24] is equivalent to

$$u_{|E} = v_{|E} : \mathcal{L}^d-\text{a.e.}$$

**Proof** See appendix.

Remark 1 Note that taking the measure-theoretic interior $(E)^1$ is of central importance. The corollary does not hold when replacing the integral over $(E)^1$ with the integral over $E$, as can be seen from the example of the closed unit ball, i.e., $E = B_1(0)$, $u = 1_E$ and $v \equiv 1$. 

Using these propositions, we now formulate the main result of this section: Alg. 1 almost surely generates an integral labeling that is of bounded variation.

**Theorem 2** Let $u \in BV(\Omega)^1$ and $f(\bar{u})$ as in Def. [1]. Then

$$\mathbb{P}(f(\bar{u}) < \infty) = 1.$$  

**Proof** The first part is to show that $(u^k)$ becomes stationary almost surely, i.e.,

$$\mathbb{P}(\exists k \in \mathbb{N} : U^k = \emptyset) = 1.$$  

Assume there exists $k$ such that $e^T c^k < 1$, and assume further that $U^k \neq \emptyset$, i.e., there exists some $x \in U^k$. Then $u^k(x) \leq c^k_j$ for all labels $j$. But then $e^T u(x) \leq e^T c^k < 1$, which is a contradiction to $u(x) \in \Delta_l$. Therefore $U^k$ must be empty. From this observation and Prop. 2, we conclude, for all $k' \in \mathbb{N}$,

$$1 \geq \mathbb{P}(\exists k \in \mathbb{N} : U^k = \emptyset) \geq \mathbb{P}(e^T c^{k'} < 1) \quad \forall k',$$

which proves (79).

In order to show that $f(\bar{u}) < \infty$ with probability 1, it remains to show that the result is almost surely in $BV(\Omega)^1$. A sufficient condition is that almost surely all iterates $u^k$ are elements of $BV(\Omega)^1$.

$$\mathbb{P}(u^k \in BV(\Omega)^1 \ \forall k \in \mathbb{N}) = 1.$$  

This is shown by Prop. 2. Then

$$\mathbb{P}(f(\bar{u}) < \infty) \geq \mathbb{P}(\{ \exists k \in \mathbb{N} : U^k = \emptyset \} \setminus \{ u^k \in BV(\Omega)^1 \ \forall k \in \mathbb{N} \})$$

$$= \mathbb{P}(u^k \in BV(\Omega)^1 \ \forall k \in \mathbb{N})$$

$$- \mathbb{P}(\{ \forall k \in \mathbb{N} : U^k \neq \emptyset \} \setminus \{ u^k \in BV(\Omega)^1 \ \forall k \in \mathbb{N} \})$$

$$= \mathbb{P}(u^k \in BV(\Omega)^1 \ \forall k \in \mathbb{N}) = 0$$

$$= 1.$$

Thus $\mathbb{P}(f(\bar{u}) < \infty) = 1$, which proves the assertion. \qed

**4 Proof of the Main Theorem**

Optimality Bounds for a Variational Relaxation of the Image Partitioning Problem
4.1 Proof of Theorem 1

In Sect. 3.2 we have shown that the rounding process induced by Alg. 1 is well-defined in the sense that it returns an integral solution \( u_* \in BV(\Omega)^l \) almost surely. We now return to proving an upper bound for the expectation of \( f(u) \) as in the approximate coarea formula \([3] \). We first show that the expectation of the linear part (data term) of \( f \) is invariant under the rounding process.

**Proposition 7** The sequence \((u^k)\) generated by Alg. 1 satisfies
\[
E(\langle u^k, s \rangle) = \langle u, s \rangle \quad \forall k \in \mathbb{N}.
\]

**Proof** In Alg. 1, instead of step 5 we consider the simpler update
\[
u^k \leftarrow e^{\gamma} 1_{\{u^{k-1}_i > \alpha^k\}} + u^{k-1} 1_{\{u^{k-1}_i \leq \alpha^k\}}.
\]

This yields exactly the same sequence \((u^k)\), since if \( u^{k-1}_i(x) > \alpha^k \) for any \( \alpha^k \geq 0 \), then either \( x \in U^{k-1} \), or \( u^{k-1}_i(x) = 1 \). In both algorithms, points that are assigned a label \( e^{\gamma} \) at some point in the process will never be assigned a different label at a later point. This is made explicit in Alg. 1 by keeping track of the set \( U^k \) of yet unassigned points. In contrast, using the step \((93)\), a point may formally be assigned the same label multiple times.

Denote \( \gamma' := (\gamma^1, \ldots, \gamma^{k-1}) \) and \( u' := u^{k-1} \). We apply induction on \( k \): For \( k \geq 1 \),
\[
E_{\gamma}(\langle u^k, s \rangle) = E_{\gamma} \frac{1}{l} \sum_{i=1}^l \int \sum_{j=1}^l s_j \cdot \left( e^{\gamma'} 1_{\{u'^j_i > \alpha^k\}} + v^{k-1} 1_{\{u'^j_i \leq \alpha^k\}} \right) dx \alpha
\]
\[
= E_{\gamma} \frac{1}{l} \sum_{i=1}^l \int \left( s_i \cdot \left( e^{\gamma'} 1_{\{u'^i_i > \alpha^k\}} + v^{k-1} 1_{\{u'^i_i \leq \alpha^k\}} \right) \langle u', s \rangle \right) dx
\]
\[
= E_{\gamma} \frac{1}{l} \sum_{i=1}^l \int \left( 1 - 1_{\{u'^i_i > \alpha^k\}} \right) \langle u', s \rangle dx \alpha.
\]

This leads to
\[
E_{\gamma}(\langle u^k, s \rangle) = E_{\gamma} \frac{1}{l} \sum_{i=1}^l \left( s_i u'^i_i + \langle u', s \rangle - u'^i_i \langle u', s \rangle \right) dx \alpha
\]
and therefore, using \( u'^i_i(x) \in \Delta_i \),
\[
E_{\gamma}(\langle u^k, s \rangle) = E_{\gamma}(\langle u^{k-1}, s \rangle).
\]
Since \( \langle u^0, s \rangle = \langle u, s \rangle \), the assertion follows by induction. \( \Box \)

**Remark 2** Prop. 7 shows that the data term is – in the mean – not affected by the probabilistic rounding process, i.e., it satisfies an exact coarea-like formula, even in the multiclass case.

Bounding the regularizer is more involved: For \( \gamma^k = (\gamma^i, \alpha^k) \), define
\[
U_{\gamma^k} := \{ x \in \Omega | u^k_i(x) \leq \alpha^k \},
\]
\[
V_{\gamma^k} := (U_{\gamma^k})^1,
\]
\[
V^k := (U^k)^1.
\]

As the measure-theoretic interior is invariant under \( \mathcal{L}^d \)-negligible modifications, given some fixed sequence \( \gamma \) the sequence \( (V^k) \) is invariant under \( \mathcal{L}^d \)-negligible modifications of \( u = u^0 \), i.e., it is uniquely defined when viewing \( u \) as an element of \( L^1(\Omega)^l \). Some calculations yield
\[
U^k = U_{\gamma^1} \cap \ldots \cap U_{\gamma^k}, \quad k \geq 1,
\]
\[
U^{k-1} \setminus U^k = U_{\gamma^2} \cap \ldots \cap U_{\gamma^{k-1}} \setminus \left( \bigcup_{i \leq 2} U_{\gamma^i} \cap \ldots \cap U_{\gamma^k} \right), \quad k \geq 2.
\]

From these observations and Prop. 4,
\[
V^k = V_{\gamma^1} \cap \ldots \cap V_{\gamma^k}, \quad k \geq 1,
\]
\[
V^{k-1} \setminus V^k = V_{\gamma^2} \cap \ldots \cap V_{\gamma^{k-1}} \setminus \left( \bigcup_{i \leq 2} V_{\gamma^i} \cap \ldots \cap V_{\gamma^k} \right), \quad k \geq 2.
\]

Moreover, since \( V^k \) is the measure-theoretic interior of \( U^k \), both sets are equal up to an \( \mathcal{L}^d \)-negligible set (cf. \([12] \)).

We now prepare for an induction argument on the expectation of the regularizing term when restricted to the sets \( V^{k-1} \setminus V^k \). The following proposition provides the initial step (\( k = 1 \)).

**Proposition 8** Assume that \( \Psi \) satisfies the lower- and upper-boundedness conditions \([23 \text{ and } 24] \). Then
\[
E \int_{V^1 \setminus V^0} d\Psi(Du) \leq \frac{2 \lambda_u}{l \lambda \Omega} \int_{\Omega} d\Psi(Du).
\]
Proof Denote \((i, \alpha) = \gamma^1\). Since \(1_{U(i, \alpha)} = 1_{V(i, \alpha)}\) \(L^d\)-a.e., we have
\[
\bar{u}_\gamma = 1_{V(i, \alpha)}e^i + 1_{D\setminus V(i, \alpha)}\bar{u}_\gamma, \quad L^d\text{- a.e.}
\] (114)
Therefore, since \(V^0 = (V^0)^1 = (\Omega)^1 = \Omega\),
\[
\int_{V^0 \cap V^1} d\Psi(D\bar{u}_\gamma) = \int_{D\setminus V(i, \alpha)} d\Psi(D\bar{u}_\gamma)
\]
\[
= \int_{D\setminus V(i, \alpha)} d\Psi\left(D\left(1_{V(i, \alpha)}e^i + 1_{D\setminus V(i, \alpha)}\bar{u}_\gamma\right)\right).
\] (115)
Since \(u \in BV(\Omega)\), we know that \(\text{Per}(V(i, \alpha)) < \infty\) holds for \(L^1\)-a.e. \(\alpha\) and any \(i\) [2 Thm. 3.40]. Therefore, we conclude from Prop. 3 that for \(L^1\)-a.e. \(\alpha\),
\[
\int_{D\setminus V(i, \alpha)} d\Psi(D\bar{u}_\gamma) \leq \lambda_u \text{ Per}(V(i, \alpha))
\]
\[
\int_{D\setminus V(i, \alpha)} d\Psi(D\bar{u}_\gamma) \leq \lambda_u \text{ Per}(V(i, \alpha)).
\] (116)
Both of the integrals are zero, since \(D\gamma^i = 0\) and \((\Omega \setminus V(i, \alpha))^0 = (V(i, \alpha))^1 = V(i, \alpha)\),
\[
\text{Carrying the bound over to the expectation yields}
\]
\[
E_\gamma \int_{D\setminus V(i, \alpha)} d\Psi(D\bar{u}_\gamma) \leq \frac{1}{l} \sum_{i=1}^{l} \int_0^1 \lambda_u \text{ Per}(V(i, \alpha)) d\alpha.
\] (117)
Also, \(\text{Per}(V(i, \alpha)) = \text{Per}(U(i, \alpha))\) since the perimeter is invariant under \(L^d\)-negligible modifications. The assertion then follows using \(V^0 = \Omega, V^1 = V(i, \alpha)\) and the coarea formula:
\[
\int_{V^0 \cap V^1} d\Psi(D\bar{u}_\gamma)
\leq \frac{1}{l} \sum_{i=1}^{l} \int_0^1 \lambda_u \text{ Per}(U(i, \alpha)) d\alpha
\] (120)
\[
\text{coarea } \frac{\lambda_u}{l} \int_{i=1}^{l} TV(u_i) = \frac{\lambda_u}{l} \int_{i=1}^{l} \|D\bar{u}_\gamma\|^2_2
\] (121)
\[
\leq 2\frac{\lambda_u}{l} \int_{\gamma} d\Psi(Du).
\] (122)
We now take care of the induction step for the regularizer bound.

**Proposition 9** Let \(\Psi\) satisfy the upper-boundedness condition (24). Then, for any \(k \geq 2\),
\[
F := E \int_{V_{k-1} \setminus V_k} d\Psi(D\bar{u})
\]
\[
\leq \frac{(l-1)}{l} E \int_{V_{k-2} \setminus V_{k-1}} d\Psi(D\bar{u}).
\] (123)
Proof Define the shifted sequence \(\gamma^k = (\gamma^k)^{\infty}_{k=1}\) by \(\gamma^k = \gamma^{k+1}\), and let
\[
W_{\gamma^k} := V_{\gamma^k} \setminus V_{\gamma^{k-1}}
\]
(125)
By Prop. 2 we may assume that, under the expectation, \(\bar{u}_\gamma\) exists and is an element of \(BV(\Omega)^k\). We denote \(\gamma^1 = (i, \alpha)\), then \(V_{\gamma^k} \setminus V_k = V(i, \alpha) \cap W_{\gamma^k}\) due to (111). For each pair \((i, \alpha)\) we denote by \((((i, \alpha), \gamma^k)\) the sequence obtained by prepending \((i, \alpha)\) to the sequence \(\gamma^k\). Then
\[
F = \frac{1}{l} \sum_{i=1}^{l} \int_0^1 \left(E_{\gamma^k} \int_{V(i, \alpha) \cap W_{\gamma^k}} d\Psi(D\bar{u}(i, \alpha), \gamma^k)\right) d\alpha.
\] (127)
Since in the first iteration of the algorithm no points in \(U(i, \alpha)\) are assigned a label, \(\bar{u}_{(i, \alpha), \gamma^k} = \bar{u}_{\gamma^k}\) holds on \(U(i, \alpha)\), and therefore \(L^1\)-a.e. on \(V(i, \alpha)\). Therefore we may apply Prop. 3 and substitute \(D\bar{u}_{(i, \alpha), \gamma^k}\) by \(D\bar{u}_{\gamma^k}\) in (127):
\[
F = \frac{1}{l} \sum_{i=1}^{l} \int_0^1 \left(E_{\gamma^k} \int_{V(i, \alpha) \cap W_{\gamma^k}} d\Psi(D\bar{u}_{\gamma^k})\right) d\alpha.
\] (128)
By definition of the measure-theoretic interior [28], the indicator function \(1_{V(i, \alpha)}\) is bounded from above by the density function \(\Theta_{U(i, \alpha)}\) of \(U(i, \alpha)\),
\[
1_{V(i, \alpha)}(x) \leq \Theta_{U(i, \alpha)}(x) := \lim_{\delta \downarrow 0} \frac{|B(x) \cap U(i, \alpha)|}{|B(x)|},
\] (130)
which exists \(H^{d-1}\)-a.e. on \(\Omega\) by [2 Prop. 3.61]. Therefore, denoting by \(B_\delta(\cdot)\) the mapping \(x \in \Omega \mapsto B_\delta(x)\),
\[
F \leq \frac{1}{l} \sum_{i=1}^{l} \int_0^1 \left(E_{\gamma^k} \int_{W_{\gamma^k} \setminus V_{\gamma^{k-1}}} \frac{|B_\delta(\cdot) \cap U(i, \alpha)|}{|B_\delta(\cdot)|} d\Psi(D\bar{u}_{\gamma^k})\right) d\alpha.
\] (131)
Rearranging the integrals and the limit, which can be justified by \(TV(\bar{u}_{\gamma^k}) < \infty\) almost surely and dominated convergence using (24), we get
\[
F \leq \frac{1}{l} \sum_{i=1}^{l} \int_0^1 \left(E_{\gamma^k} \lim_{\delta \downarrow 0} \int_{W_{\gamma^k}} \frac{|B_\delta(\cdot) \cap U(i, \alpha)|}{|B_\delta(\cdot)|} d\Psi(D\bar{u}_{\gamma^k})\right) d\alpha
\]
\[
= \frac{1}{l} \sum_{i=1}^{l} \int_0^1 \left(E_{\gamma^k} \lim_{\delta \downarrow 0} \int_{W_{\gamma^k}} \frac{1}{|B_\delta(\cdot)|} \right) \left(\int_{1_{(u_i(\alpha) \leq \alpha)\} dy d\alpha}\right) d\Psi(D\bar{u}_{\gamma^k}).
\] (131)
We again apply Proposition 1.78 to the two innermost integrals (alternatively, use Fubini’s theorem), which leads to

\[
F \leq \frac{1}{l} \mathbb{E}_\gamma \lim_{\delta \to 0} \int_{W_\gamma} \frac{1}{|\mathcal{B}_\delta(\cdot)|} \left( \sum_{i=1}^l \int_{\mathcal{B}_\delta(y)} (1 - u_i(y))dy \right) d\Psi(D\bar{u}_\gamma). \tag{132}
\]

Using the fact that \( u(y) \in \mathcal{D}_t \), this collapses according to

\[
F \leq \frac{1}{l} \mathbb{E}_\gamma \lim_{\delta \to 0} \int_{W_\gamma} \frac{1}{|\mathcal{B}_\delta(\cdot)|} \left( \int_{\mathcal{B}_\delta(y)} (l - 1)dy \right) d\Psi(D\bar{u}_\gamma) \tag{134}
= \frac{l - 1}{l} \mathbb{E}_\gamma \int_{W_\gamma} d\Psi(D\bar{u}_\gamma) \tag{135}
= \frac{l - 1}{l} \mathbb{E}_\gamma \int_{V_{k-2} \setminus V_{k-1}} d\Psi(D\bar{u}_\gamma). \tag{136}
\]

Reverting the index shift and using \( \bar{u}_\gamma = \bar{u} \) concludes the proof:

\[
F \leq \frac{l - 1}{l} \mathbb{E}_\gamma \int_{V_{k-1} \setminus V_k} d\Psi(D\bar{u}_\gamma). \tag{137}
\]

We are now ready to prove the main result, Theorem 1, as stated in the introduction.

**Proof (Theorem 1)** The fact that the algorithm provides \( \bar{u} \in \mathcal{G}_\gamma \) almost surely follows from Theorem 2. Therefore there almost surely exists \( k' := k'(\gamma) \geq 1 \) such that \( U^{k'} = 0 \) and \( \bar{u}_\gamma = u^{k'} \). On one hand, this implies

\[
\langle u^{k'}, s \rangle = \lim_{k \to \infty} \langle u^{k'}, s \rangle \tag{138}
\]

almost surely. On the other hand, \( V^{k'} = (U^{k'})^1 = 0 \) and therefore

\[
\bigcup_{k=1}^{k'} V^{k-1} \setminus V^k = \Omega \setminus V^{k'} = \Omega \tag{139}
\]

almost surely. The equality (*) can be shown by induction: For the base case \( k' = 1 \), we have \( V^0 = (U^0)^1 = (\Omega)^1 = \Omega \), since \( \Omega \) is the open unit box. For \( k' \geq 2 \),

\[
\bigcup_{k=1}^{k'} V^{k-1} \setminus V^k = \bigcup_{k=1}^{k'} (V^{k-1} \setminus V^{k-1}) \tag{140}
= \bigcup_{k=1}^{k'} (V^{k-1} \setminus V^k) \tag{141}
= \bigcup_{k=1}^{k'} (V^{k-1} \setminus V^k) \cup (\Omega \setminus V^k) \tag{142}
= \Omega \setminus V^{k-1}. \tag{143}
\]

almost surely (cf. (122)). From (138) and (139) we obtain

\[
\mathbb{E}_\gamma f(\bar{u}_\gamma) = \mathbb{E}_\gamma \left( \lim_{k \to \infty} \langle u^k, s \rangle \right) + \mathbb{E}_\gamma \left( \sum_{k=1}^{\infty} \int_{V_{k-1} \setminus V_k} d\Psi(D\bar{u}_\gamma) \right) \tag{144}
= \lim_{k \to \infty} \mathbb{E}_\gamma \left( \langle u^k, s \rangle \right) + \sum_{k=1}^{\infty} \mathbb{E}_\gamma \int_{V_{k-1} \setminus V_k} d\Psi(D\bar{u}_\gamma). \tag{145}
\]

The first term in (145) is equal to \( \langle u, s \rangle \) due to Proposition 3. An induction argument using Proposition 8 and Proposition 9 shows that the second term can be bounded according to

\[
\sum_{k=1}^{\infty} \mathbb{E}_\gamma \int_{V_{k-1} \setminus V_k} d\Psi(D\bar{u}_\gamma) \leq \sum_{k=1}^{\infty} \left( (l - 1) \right) k - 2 \frac{\lambda_u}{\lambda_t} \int_\Omega d\Psi(Du) \leq 2 \frac{\lambda_u}{\lambda_t} \int_\Omega d\Psi(Du), \tag{147}
\]

therefore

\[
\mathbb{E}_\gamma f(\bar{u}_\gamma) \leq \langle u, s \rangle + 2 \frac{\lambda_u}{\lambda_t} \int_\Omega d\Psi(Du). \tag{149}
\]

Since \( s \geq 0 \) and \( \lambda_u \geq \lambda_t \), and therefore the linear term is bounded by \( \langle u, s \rangle \leq 2(\lambda_u / \lambda_t)\langle u, s \rangle \), this proves the assertion. Swapping the integral and limit in (145) can be justified retrospectively by the dominated convergence theorem, using \( 0 \leq \langle u, s \rangle \leq \infty \) and \( \int_\Omega d\Psi(Du) < \infty \) due to the upper-boundedness condition and Proposition 3.

\( \square \)

Corollary 1 (see introduction) follows immediately from \( f(u^*) \leq f(u^\gamma) \), cf. (144). We have demonstrated that the proposed approach allows to recover, from the solution \( u^* \) of the convex relaxed problem \( 13 \), an approximate integral solution \( \bar{u}^* \) of the nonconvex original problem \( 1 \) with an upper bound on the objective.

For the specific case \( \Psi = \Psi_d \), we have

**Proposition 10** Let \( d : \mathcal{I}^2 \to \mathbb{R}_{\geq 0} \) be a metric and \( \Psi = \Psi_d \). Then one may set

\[
\lambda_u = \max_{i,j \in \{1, \ldots, l\}} d(i, j) \text{ and } \lambda_t = \min_{i \neq j} d(i, j).
\]

**Proof** From the remarks in the introduction we obtain (cf. (19))

\[
\Psi_d(\nu(e^i - e^j)) = d(i, j),
\]

which shows the upper bound. For the lower bound, set \( c := \min_{i \neq j} d(i, j) \), \( v^\nu := \frac{v}{\|v\|_2} \) and \( v := v(I - \frac{1}{l} ee^\top) \).
Then \( v \in D_{\text{loc}}^d \) since \( \|v^i - v^j\|_2 = \|v'^i - v'^j\|_2 \leq c \) and \( v c = v'(I - \frac{1}{2}v v^\top)c = 0. \) Therefore, for \( w \in \mathbb{R}^d \) satisfying \( w e = 0, \)
\[
\Psi_d(w) \geq \langle w, v \rangle = \langle w', v' \rangle = \sum_{i=1}^t \langle w^i, \frac{c}{2} \frac{w^i}{\|w^i\|_2} \rangle = \frac{c}{2} \sum_{i=1}^t \|w^i\|_2,
\]
(150) proving the lower bound.

Finally, for \( \Psi_d \) we obtain the factor
\[
2 \frac{\lambda_u}{\lambda_l} = 2 \frac{\max_{i,j} d(i,j)}{\min_{i\neq j} d(i,j)},
\]
(152) determining the optimality bound, as claimed in the introduction [27]. The bound in [27] is the same as the known bounds for finite-dimensional metric labeling [12] and \( \alpha \)-expansion [4], however it extends these results to problems on continuous domains for a broad class of regularizers.

5 Conclusion

In this work we considered a method for recovering approximate solutions of image partitioning problems from solutions of a convex relaxation. We proposed a probabilistic rounding method motivated by the finite-dimensional framework, and showed that it is possible to obtain \textit{a priori} bounds on the optimality of the integral solution obtained by rounding a solution of the convex relaxation.

The obtained bounds are compatible with known bounds for the finite-dimensional setting. However, to our knowledge, this is the first fully convex approach that is both formulated in the spatially continuous setting and provides a true \textit{a priori} bound. We showed that the approach can also be interpreted as an approximate variant of the coarea formula.

While the results apply to a quite general class of regularizers, they are formulated for the homogeneous case. Non-homogeneous regularizers constitute an interesting direction for future work. In particular, such regularizers naturally occur when applying convex relaxation techniques [1][25] in order to solve nonconvex variational problems.

With the increasing computational power, such techniques have become quite popular recently. For problems where the convexity is confined to the data term, they permit to find a global minimizer. A proper extension of the results outlined in this work may provide a way to find good approximate solutions of problems where also the \textit{regularizer} is nonconvex.

6 Appendix

**Proof (Prop. 3)** In order to prove the first assertion [27], note that the mapping \( w \mapsto \Psi(w^+) \) is convex, therefore it must assume its maximum on the polytope \( \Delta_l - \Delta_l := \{ z^1 - z^2, z^2 - \Delta_l \} \) in a vertex of the polytope. Since the polytope \( \Delta_l - \Delta_l \) is the difference of two polytopes, its vertex set is at most the difference of their vertex sets, \( V := \{ e^i - e^j | i, j \in \{1, \ldots, l\} \}. \) On this set, the bound \( \Psi(w^+) \leq \lambda_u \) holds for \( w \in V \) due to the upper-boundedness condition (24), which proves (57).

The second equality (89) follows from the fact that \( G := \{ b^{ik} := e^k(e^i - e^{i+1})^\top | 1 \leq k \leq d, 1 \leq i \leq l - 1 \} \) is a basis of the linear subspace \( W, \) satisfying \( \Psi(b^{ik}) \leq \lambda_u, \) and \( \Psi \) is positively homogeneous and convex, and thus subadditive. Specifically, there is a linear transform \( T : W \rightarrow \mathbb{R}^{x(l-1)} \) such that \( w = \sum_{i,k} b^{ik} \alpha_{ik} \) for \( \alpha = T(w). \) Then
\[
\Psi(w) = \Psi \left( \sum_{i,k} b^{ik} \alpha_{ik} \right) = \leq \Psi \left( \sum_{i,k} |\alpha_{ik}| \text{sgn}(\alpha_{ik}) b^{ik} \right) \leq \sum_{i,k} |\alpha_{ik}| \Psi \left( \text{sgn}(\alpha_{ik}) b^{ik} \right).
\]
(154) Since (24) provides \( \Psi(\pm b^{ik}) \leq \lambda_u, \) we obtain
\[
\Psi(w) \leq \lambda_u \sum_{i,k} |\alpha_{ik}| \leq \lambda_u \|T\|\|w\|_2
\]
(156) for some suitable operator norm \( \| \cdot \| \) and any \( w \in W. \)

**Proof (Prop. 4)** We prove mutual inclusion:
\( \alpha \subseteq \alpha' \): From the definition of the measure-theoretic interior,
\[
x \in (E \cap F)^1 \Rightarrow \lim_{\delta \searrow 0} \frac{|B_{\delta}(x) \cap E \cap F|}{|B_{\delta}(x)|} = 1.
\]
(157) Since \( |B_{\delta}(x)| \geq |B_{\delta}(x) \cap E| \geq |B_{\delta}(x) \cap E \cap F| \) (and vice versa for \( |B_{\delta}(x) \cap F| \)), it follows by the “sandwich” criterion that both \( \lim_{\delta \searrow 0} |B_{\delta}(x) \cap E|/|B_{\delta}(x)| \) and \( \lim_{\delta \searrow 0} |B_{\delta}(x) \cap F|/|B_{\delta}(x)| \) exist and are equal to 1, which shows \( x \in E^1 \cap F^1. \)
\( \alpha' \supseteq \alpha \): Assume that \( x \in E^1 \cap F^1. \) Then
\[
1 \geq \limsup_{\delta \searrow 0} \frac{|B_{\delta}(x) \cap E \cap F|}{|B_{\delta}(x)|} \geq \liminf_{\delta \searrow 0} \frac{|B_{\delta}(x) \cap E \cap F|}{|B_{\delta}(x)|} \geq \liminf_{\delta \searrow 0} \frac{|B_{\delta}(x) \cap E| + |B_{\delta}(x) \cap F| - |B_{\delta}(x) \cap (E \cup F)|}{|B_{\delta}(x)|} \geq 1,
\]
(158) (159)
We obtain equality,

\[
1 \geq \lim_{\delta \searrow 0} \inf \frac{|B_\delta(x) \cap E \cap F|}{|B_\delta(x)|} \geq \lim_{\delta \searrow 0} \frac{|B_\delta(x) \cap E|}{|B_\delta(x)|} + \lim_{\delta \searrow 0} \frac{|B_\delta(x) \cap F|}{|B_\delta(x)|} + \lim_{\delta \searrow 0} \frac{\left(-\frac{|B_\delta \cap (E \cup F)|}{|B_\delta(x)|}\right)}{\geq 1,}
\]

from which we conclude that

\[
\lim_{\delta \searrow 0} \sup \frac{|B_\delta(x) \cap E \cap F|}{|B_\delta(x)|} = \lim_{\delta \searrow 0} \frac{|B_\delta(x) \cap E \cap F|}{|B_\delta(x)|} = 1,
\]

i.e., \(x \in (E \cap F)^c\).

**Proof (Prop. 3)** First note that

\[
\int_{E \cap \Omega} \|w_{E \cap \Omega}^x - w_{E \cap \Omega}^y\| d\mathcal{H}^{d-1} \leq \sup \|w_{E \cap \Omega}^x - w_{E \cap \Omega}^y\| \cdot |E \cap \Omega|^{d-1} (\mathcal{H}^{d-1}(E \cap \Omega)),
\]

\[
(*) \leq \sqrt{2} \text{TV}(1_E) < \infty.
\]

The inequality \((*)\) is a consequence of the definition of \(w_{E \cap \Omega}^x\) and Thm. 3.59, and \((**)\) follows directly from \(w(x), w(y) \in \Delta_1\). The upper bound permits applying Thm. 3.84 on \(w\), which provides \(w \in \text{BV}(\Omega)^1\) and [2]. Due to Prop. 3.61, the sets \((E)^0, (E)^1\) and \(E \cap \Omega\) form a (pairwise disjoint) partition of \(\Omega\), up to an \(\mathcal{H}^{d-1}\)-zero set. Moreover, since \(\Psi(Du) \leq |Du| \leq \mathcal{H}^{d-1}\) by construction, we have, for some Borel set \(A\),

\[
\int_A \Psi(Dw) = \int_{A \cap (E)^1} d\Psi(Dw) + \int_{A \cap (E)^0} d\Psi(Dw) + \int_{A \cap E \cap \Omega} \left(\nu_E (w^x_{E \cap \Omega}(x) - w^y_{E \cap \Omega}(x)) \right) d\mathcal{H}^{d-1} \leq \int_{A \cap (E)^1} d\Psi(Dw) + \int_{A \cap (E)^0} d\Psi(Dw) + \int_{A \cap E \cap \Omega} \lambda_a d\mathcal{H}^{d-1} \leq 
\]

The inequality \((**)\) holds due to the upper boundedness and Prop. 3. From [2] Prop. 2.37] we obtain that \(\Psi\) is additive on mutually singular Radon measures \(\mu, \nu\), i.e., if \(|\mu|, |\nu|\), then

\[
\int_B d\Psi(\mu + \nu) = \int_B d\Psi(\mu) + \int_B d\Psi(\nu)
\]

for any Borel set \(B \subseteq \Omega\). Substituting \(Du\) in [171] according to [172] and using the fact that the three measures that form \(Du\) in [172] are mutually singular, the additivity property [172] leads to the remaining assertion,

\[
\int_A d\Psi(Dw) \leq \int_{A \cap (E)^1} d\Psi(Du) + \int_{A \cap (E)^0} d\Psi(Dv) + \lambda_a \text{Per}(E).
\]

**Proof (Prop. 4)** We first show [90]. It suffices to show that

\[
\{x \in (E)^1 \iff x \in E\} \text{ for } \mathcal{L}^d\text{-a.e. } x \in \Omega.
\]

This can be seen by considering the precise representative \(1_E\) of \(1_E\) [2] Def. 3.63: Starting with the definition,

\[
x \in (E)^1 \iff \lim_{\delta \searrow 0} \frac{|E \cap B_\delta(x)|}{|B_\delta(x)|} = 1,
\]

the fact that \(\lim_{\delta \searrow 0} \frac{|\Omega \cap B_\delta(x)|}{|B_\delta(x)|} = 1\) implies

\[
x \in (E)^1 \iff \lim_{\delta \searrow 0} \frac{|\Omega \cap B_\delta(x)|}{|B_\delta(x)|} = 0
\]

Substituting \(E\) by \(\Omega \cap E\), the same equivalence shows that \(x \in (E)^0 \iff 1_{\Omega \cap E}(x) = 1 \iff 1_E(x) = 0\). As \(\mathcal{L}^d(\Omega \setminus ((E)^0 \cup (E)^1)) = 0\), this shows that \(1_E = 1_E\) \(\mathcal{L}^d\text{-a.e.}\). Using the fact that \(1_E = 1_E\) [2] Prop. 3.64], we conclude that \(1_{(E)^1} = 1_E\) \(\mathcal{L}^d\text{-a.e.}\), which proves [174] and therefore the assertion [90].
e.g., Prop. 3.2). We continue from (181) via
\[ D_u(v)(E) \]
and
\[ \nu_v(t) = \frac{1}{t} \wedge \varphi(t) \]
(182)

\[ \varphi(t) = 1 - \left( 1 - \frac{1}{t} \right) \wedge \varphi(t) \]
(183)

\[ \quad \nu_v(t) = \frac{1}{t} \wedge \varphi(t) \]
(184)

\[ \Psi(D_u)(E) = \Psi(D_u)(E) \]
(185)

\[ \Psi(D_u)(E) = \Psi(D_u)(E) \]
(186)

Therefore \( D_u(v)(E) = D_v(v)(E) \). Then,
\[ \Psi(D_u)(E) = \Psi(D_u)(E) \]
(187)

\[ \Psi(D_u)(E) = \Psi(D_u)(E) \]
(188)

In the equality (*) we used the additivity of \( \Psi \) on mutually singular Radon measures [2, Prop. 2.37]. By definition of the total variation, \( |\mu| = |\mu| \) holds for any measure \( \mu \), therefore \( |Du|(\Omega \setminus E) = |Du|(\Omega \setminus E) \) and \( |Du|(\Omega \setminus E) = 0 \), which together with again by definition \( |\mu| \leq |\mu| \) implies that the second term in (185) vanishes. Since all observations equally hold for \( v \) instead of \( u \), we conclude
\[ \Psi(D_u)(E) = \Psi(D_u)(E) \]
(189)

\[ \Psi(D_u)(E) = \Psi(D_u)(E) \]
(190)

\[ \Psi(D_u)(E) = \Psi(D_u)(E) \]
(191)

Equation (185) follows immediately.

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