

Learning Belief Functions from Data: Experimental Supplement

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Introduction

This notebook provides a compact computational companion to the paper

“Learning Belief Functions from Data via Polyhedral Methods” (Kybernetika, 2025).

Its purpose is to **reproduce and document all worked examples from the paper**:

- construction of **Jeffreys-based pseudo-belief functions** g from count data,
- application of the **Upper Approximation Procedure (UAP)** to these functions (Examples 1–3, Tables 1, 4–6),
- and the **polyhedral LP approach** used to select dominating belief functions (Examples 4–5, Tables 7–9).

In particular, the notebook:

1. Shows how Jeffreys’ intervals turn simple frequency tables into a normalized, monotone set function g on $\mathcal{P}(\Omega)$.
2. Applies the Upper Approximation Procedure to g and its modifications, exactly reproducing Tables 1, 4, 5, and 6.
3. Constructs the dominance polytope \mathcal{M}_g and uses several linear objectives (L1, HD, CW) to obtain LP-based upper approximations, reproducing Tables 7–9.
4. For the **quaternary Example 5**, compares the LP solutions with the outcome of UAP in a single summary table.

The notebook therefore serves as an executable “log file” for all numerical tables in the paper and as a reference implementation of the UAP and LP-based upper approximation methods.

```
# Required packages
library(lpSolve)
library(knitr)

# =====
# Auxiliary functions
# =====

# Generate all subsets
powerset <- function(Omega) {
  n <- length(Omega)
  unlist(lapply(0:n, function(i) combn(Omega, i, simplify = FALSE)), recursive = FALSE)
}

subset_name <- function(S) {
  if (length(S)==0) return("{}")
  paste0("{", paste(S, collapse=","), "}")
}
```

Jeffreys Lower Bounds

To construct the statistical lower approximation $g(A)$ on the Boolean algebra $\mathcal{P}(\Omega)$, we follow the Jeffreys-based procedure described in Section 2 of the paper.

Assume that observations are recorded on subsets $B \subseteq \Omega$, possibly ambiguous (e.g. $\{w_1, w_2\}$). Let n_B denote the number of observations associated with B , and let

$$N = \sum_{B \subseteq \Omega} n_B$$

be the total number of observed cases.

An observation assigned to B supports the event A only when $B \subseteq A$. The effective number of “successes” for event A is therefore

$$s(A) = \sum_{B \subseteq A} n_B.$$

Under Jeffreys’ prior $\text{Beta}(1/2, 1/2)$, the posterior distribution of the binomial parameter $p(A)$ is

$$p(A) \mid \text{data} \sim \text{Beta}\left(s(A) + \frac{1}{2}, N - s(A) + \frac{1}{2}\right),$$

and the corresponding lower $(1 - \alpha)$ -credible bound is

$$g(A) = L(A) = \text{qbeta}\left(\frac{\alpha}{2}, s(A) + \frac{1}{2}, N - s(A) + \frac{1}{2}\right).$$

Following the article, we additionally impose the normalization condition $g(\Omega) = 1$, which ensures that the resulting set function behaves as a *pseudo-belief function* compatible with the dominance relation used later in the Upper Approximation Procedure.

All computations of $g(A)$ in this notebook use the function `jeffreys_pseudo_belief_function()` defined below, which implements the mapping $\{n_B\} \mapsto g(\cdot)$ exactly as described above.

For convenience, we include the empty set \emptyset in the internal representation (with $g(\emptyset) = 0$, but we omit it from the printed tables.

```
# -----
# Jeffreys confidence interval for the binomial model
# Returns the lower and upper (1 - alpha) credible bounds.
# -----
jeffreys_interval <- function(successes, total, alpha = 0.05,
                             a0 = 0.5, b0 = 0.5) {

  lower <- qbeta(alpha / 2, successes + a0, total - successes + b0)
  upper <- qbeta(1 - alpha / 2, successes + a0, total - successes + b0)

  list(lower = lower, upper = upper)
}

# -----
# Normalize a named vector of counts:
# Ensures all subsets exist; missing subsets are assigned zero.
# -----
normalize_counts <- function(counts_named, subsets) {

  subset_names <- sapply(subsets, subset_name)
  out <- setNames(rep(0, length(subsets)), subset_names)

  if (is.null(names(counts_named)))
    stop("counts_named must be a named vector, e.g. c('{w1}'=5, '{w1,w2}'=2).")

  # fill matching names
  out[names(counts_named)] <- as.numeric(counts_named)

  out
}

# -----
# Construct Jeffreys-based pseudo-belief function g(A):
# g(A) = Jeffreys lower bound for f(A)
# where "successes" for A is the total number of observations
# assigned to subsets B \subseteq A.
#
# The function automatically enforces g(Omega) = 1, consistently with the paper
# -----
jeffreys_pseudo_belief_function <- function(Omega, counts_named, alpha = 0.05) {
```

```

# all subsets in cardinality order
subsets <- powerset(Omega)
subset_names <- sapply(subsets, subset_name)

# full count vector for all subsets
counts_full <- normalize_counts(counts_named, subsets)
N <- sum(counts_full)

# compute "successes"  $s(A) = \sum_{B \setminus \text{subteq } A} n_B$ 
successes <- sapply(subsets, function(A) {
  idx <- sapply(subsets, function(B) all(B %in% A))
  sum(counts_full[idx])
})

# allocate vector for  $g(A)$ 
g_values <- numeric(length(subsets))

for (i in seq_along(subsets)) {
  A <- subsets[[i]]

  # by convention  $g(\text{emptyset}) = 0$ 
  if (length(A) == 0) {
    g_values[i] <- 0
  } else {
    ji <- jeffreys_interval(successes[i], N, alpha = alpha)
    g_values[i] <- ji$lower
  }
}

# normalize:  $g(\text{Omega})$  must equal 1
g_values[length(g_values)] <- 1

list(
  g          = g_values,
  subsets    = subsets,
  subset_names = subset_names,
  counts     = counts_full,
  N          = N
)
}

```

Example: Ambiguous observations

This example illustrates how the Jeffreys-based construction processes **partially ambiguous measurements**.

We consider 30 observations in total:

- 25 observations are fully specified (assigned to the singletons $\{w_1\}, \{w_2\}, \{w_3\}$),
- 3 observations cannot distinguish between states w_1 and w_2 and are therefore recorded as the set $\{w_1, w_2\}$,
- 2 observations provide no discriminatory information at all, represented by the full set $\{w_1, w_2, w_3\}$.

In this setting, an ambiguous observation recorded as B contributes evidence to every event A such that $B \subseteq A$.

Consequently, the resulting Jeffreys lower bounds $g(A)$ naturally incorporate both precise and imprecise

information without forcing an artificial resolution of uncertainty.

```
## Frame of discernment
Omega <- c("w1", "w2", "w3")

## Example: 30 observations (25 precise, 3 ambiguous {w1,w2}, 2 fully ambiguous)
counts <- c(
  "{w1}"      = 10,
  "{w2}"      = 8,
  "{w3}"      = 7,
  "{w1,w2}"   = 3,
  "{w1,w2,w3}" = 2
)

## Compute Jeffreys pseudo-belief function g(A)
g_result <- jeffreys_pseudo_belief_function(Omega, counts, alpha = 0.05)

## Create output table (drop empty set)
df <- data.frame(
  subset = g_result$subset_names,
  g       = round(g_result$g, 3),
  row.names = NULL
)

df <- df[df$subset != "{}", ] # hide the empty set

df
```

##	subset	g
## 2	{w1}	0.186
## 3	{w2}	0.135
## 4	{w3}	0.111
## 5	{w1,w2}	0.523
## 6	{w1,w3}	0.390
## 7	{w2,w3}	0.328
## 8	{w1,w2,w3}	1.000

Upper Approximation Procedure (UAP)

To obtain a belief function f that dominates the statistically derived pseudo-belief function g , the paper employs the **Upper Approximation Procedure** (UAP) introduced in Section 3.

The method constructs a mass assignment m and the corresponding belief function f by processing the subsets of Ω in increasing order of cardinality.

For every set A , the value of $f(A)$ is determined so that

$$f(A) \geq g(A) \quad \text{and} \quad f(A) \geq \sum_{B \subset A} m(B),$$

where the second inequality ensures the internal consistency of the belief function. The increment

$$m(A) = f(A) - \sum_{B \subset A} m(B)$$

defines the basic probability assigned to A .

If the procedure reaches the top element Ω and produces $f(\Omega) > 1$, the construction fails; otherwise the output is a valid belief function dominating g .

This behaviour matches the examples in the article, including the failure for Table 4 and the successful runs for Table 1 (modified), Table 5, and Table 6.

All computations below use the implementation `upper_approximation()`, which follows the pseudocode of the article exactly. The procedure processes subsets in increasing cardinality, ensuring the Möbius recursion is respected.

```
# -----
# Upper Approximation Procedure (UAP)
#
# Input:
#   Omega      ... vector of frame elements
#   g_values    ... numeric vector of g(A) in cardinality order,
#                 or named by subset_name().
#
# Output:
#   list(success, subsets, f, m)
#
# The procedure assumes that g(Omega) = 1 (enforced here).
# -----
upper_approximation <- function(Omega, g_values) {

  # 1) Generate subsets in canonical order (cardinality-first)
  subsets <- powerset(Omega)
  subsets <- subsets[order(sapply(subsets, length))]
  subset_names <- sapply(subsets, subset_name)

  # 2) If g_values have names -> reorder; if not -> assume correct order
  if (!is.null(names(g_values))) {
    g_values <- g_values[subset_names]
  } else {
    names(g_values) <- subset_names
  }

  # 3) Enforce normalization: g(Omega) = 1
  # Find index of \Omega = full set
  full_idx <- which(subset_names == subset_name(Omega))
  g_values[full_idx] <- 1

  k <- length(subsets)
  f <- numeric(k)
  m <- numeric(k)

  # 4) Main UAP loop (exact definition from article)
  for (i in seq_len(k)) {
    A <- subsets[[i]]

    # proper subsets of A
    idx <- sapply(subsets, function(B)
      length(B) < length(A) && all(B %in% A))

    sum_prev <- sum(m[idx])
```

```

    # f(A)
    f[i] <- max(g_values[i], sum_prev)

    # m(A)
    m[i] <- f[i] - sum_prev
  }

# 5) Check failure condition
if (f[k] > 1 + 1e-12) {
  return(list(
    success = FALSE,
    message = "UAP failed: f(Omega) > 1",
    subsets = subset_names,
    f = f,
    m = m
  ))
}

# IMPORTANT:
# f(Omega) is NOT overwritten.
# m(Omega) stays as computed.

list(
  success = TRUE,
  subsets = subset_names,
  f = f,
  m = m
)
}

# =====
# Helper: Print UAP table exactly like in paper
# =====

uap_table <- function(result, g_values) {
  df <- data.frame(
    A      = result$subsets,
    g      = g_values,
    sum_prev = round(result$f - result$m, 6),
    f      = round(result$f, 6),
    m_f    = round(result$m, 6),
    row.names = NULL
  )

  # remove the first row (empty set)
  df <- df[-1, ]

  df
}

```

Table 1 (ternary, original g)

```
Omega3 <- c("w1", "w2", "w3")

g_ternary <- c(
  "{}"=0,
  "{w1}"=0.1, "{w2}"=0.1, "{w3}"=0.2,
  "{w1,w2}"=0.5, "{w1,w3}"=0.4, "{w2,w3}"=0.7,
  "{w1,w2,w3}"=1.0
)

res <- upper_approximation(Omega3, g_ternary)
res$success # expected: FALSE (as in Table 1)

## [1] FALSE

table <- uap_table(res, g_ternary)

# print table
knitr::kable(
  table,
  caption = "Table 1 - Upper Approximation for the original pseudo-belief function g on a ternary frame",
  align = "lrrrr",
  row.names = FALSE
)
```

Table 1: Table 1 — Upper Approximation for the original pseudo-belief function g on a ternary frame.

A	g	sum_prev	f	m_f
{w1}	0.1	0.0	0.1	0.1
{w2}	0.1	0.0	0.1	0.1
{w3}	0.2	0.0	0.2	0.2
{w1,w2}	0.5	0.2	0.5	0.3
{w1,w3}	0.4	0.3	0.4	0.1
{w2,w3}	0.7	0.3	0.7	0.4
{w1,w2,w3}	1.0	1.2	1.2	0.0

For the original pseudo-belief function g , the UAP fails because the lower bounds on several subsets jointly force the final value $f(\Omega)$ to exceed 1.

Table 1 (modified \tilde{g})

```
g_ternary_tilde <- c(
  "{}"=0,
  "{w1}"=0.2, "{w2}"=0.2, "{w3}"=0.3,
  "{w1,w2}"=0.5, "{w1,w3}"=0.4, "{w2,w3}"=0.7,
  "{w1,w2,w3}"=1.0
)

res_tilde <- upper_approximation(Omega3, g_ternary_tilde)
res_tilde$success # expected: TRUE

## [1] TRUE
```



```

table <- uap_table(res_tilde, g_ternary_tilde)

# print table
knitr::kable(
  table,
  caption = "Table 1 (modified) - UAP succeeds for the adjusted pseudo-belief.",
  align = "lrrrr",
  row.names = FALSE
)

```

Table 2: Table 1 (modified) — UAP succeeds for the adjusted pseudo-belief.

A	g	sum_prev	f	m_f
{w1}	0.2	0.0	0.2	0.2
{w2}	0.2	0.0	0.2	0.2
{w3}	0.3	0.0	0.3	0.3
{w1,w2}	0.5	0.4	0.5	0.1
{w1,w3}	0.4	0.5	0.5	0.0
{w2,w3}	0.7	0.5	0.7	0.2
{w1,w2,w3}	1.0	1.0	1.0	0.0

The modified function \tilde{g} , which only increases singleton values, removes this conflict: the accumulated mass from proper subsets stays below 1, allowing the UAP to construct a valid dominating belief function. The example illustrates that the feasibility of UAP is strongly dependent on the internal coherence of the statistical lower bounds.

UAP Reproduction of Tables 4, 5, 6

In the quaternary case $|\Omega| = 4$, the data originate from a simple frequency table (Table 2 in the article). Applying the Jeffreys construction yields a pseudo-belief function g whose properties depend sensitively on the confidence level α .

This section reproduces the behaviour of the Upper Approximation Procedure under three different conditions corresponding to Tables 4, 5, and 6.

```

Omega4 <- c("w1", "w2", "w3", "w4")

counts_raw <- c(
  "{w1}"=14, "{w2}"=17, "{w3}"=15, "{w4}"=6
)

```

Table 4 — Jeffreys lower bounds with $\alpha = 0.05$

For the standard 95% Jeffreys interval, many lower bounds become relatively tight, especially on two- and three-element subsets.

When fed into the UAP, the cumulative mass assigned to proper subsets grows too large, and the procedure attempts to set

$$f(\Omega) > 1,$$

which violates the normalization constraint.

Consequently, **UAP fails** for this choice of α , matching the result reported in Table 4 of the article.

This failure illustrates that the lower statistical bounds may be too optimistic for a belief function to dominate them.

```
# Compute pseudo-belief function g(A)
g4_res <- jeffreys_pseudo_belief_function(Omega4, counts_raw, alpha = 0.05)
g4_vals <- g4_res$g
names(g4_vals) <- g4_res$subset_names

# Apply UAP
res4 <- upper_approximation(Omega4, g4_vals)

# Expected: FALSE (matches article Table 4)
res4$success

## [1] FALSE

# Construct table
tab4 <- uap_table(res4, g4_vals)

# print table
knitr::kable(
  tab4,
  caption = "Table 4 - UAP failure for Jeffreys lower bounds with alpha = 0.05.",
  align = "lrrrr",
  row.names = FALSE
)
```

Table 3: Table 4 — UAP failure for Jeffreys lower bounds with $\alpha = 0.05$.

A	g	sum_prev	f	m_f
{w1}	0.1634559	0.000000	0.163456	0.163456
{w2}	0.2114491	0.000000	0.211449	0.211449
{w3}	0.1792032	0.000000	0.179203	0.179203
{w4}	0.0496248	0.000000	0.049625	0.049625
{w1,w2}	0.4605750	0.374905	0.460575	0.085670
{w1,w3}	0.4226370	0.342659	0.422637	0.079978
{w1,w4}	0.2615272	0.213081	0.261527	0.048447
{w2,w3}	0.4798459	0.390652	0.479846	0.089194
{w2,w4}	0.3134786	0.261074	0.313479	0.052405
{w3,w4}	0.2786434	0.228828	0.278643	0.049815
{w1,w2,w3}	0.7775591	0.808950	0.808950	0.000000
{w1,w2,w4}	0.5794675	0.611051	0.611051	0.000000
{w1,w3,w4}	0.5389325	0.570524	0.570524	0.000000
{w2,w3,w4}	0.6001125	0.631691	0.631691	0.000000
{w1,w2,w3,w4}	1.0000000	1.009241	1.009241	0.000000

Table 5 — Jeffreys lower bounds with $\alpha = 0.02$

Reducing the confidence level to $\alpha = 0.02$ produces wider Jeffreys intervals and therefore **smaller lower bounds** $g(A)$.

These relaxed bounds reduce the dominance pressure on the UAP recursion, allowing it to satisfy all inequalities while keeping

$$f(\Omega) = 1.$$

The UAP therefore **succeeds**, giving the belief function listed in Table 5. This demonstrates how the statistical confidence level directly influences the feasibility of constructing a dominating belief function.

```
## Table 5 - Jeffreys lower bounds (alpha = 0.02)

# Compute pseudo-belief function g(A)
g5_res <- jeffreys_pseudo_belief_function(Omega4, counts_raw, alpha = 0.02)
g5_vals <- g5_res$g
names(g5_vals) <- g5_res$subset_names

# Apply UAP
res5 <- upper_approximation(Omega4, g5_vals)

# Expected: TRUE (matches article Table 5)
res5$success

## [1] TRUE

# Construct table
tab5 <- uap_table(res5, g5_vals)

# Print table

# Knit-friendly table
knitr::kable(
  tab5,
  caption = "Table 5 - UAP succeeds for Jeffreys lower bounds with alpha = 0.02.",
  align = "lrrrr",
  row.names = FALSE
)
```

Table 4: Table 5 — UAP succeeds for Jeffreys lower bounds with alpha = 0.02.

A	g	sum_prev	f	m_f
{w1}	0.1465951	0.000000	0.146595	0.146595
{w2}	0.1923940	0.000000	0.192394	0.192394
{w3}	0.1615697	0.000000	0.161570	0.161570
{w4}	0.0408735	0.000000	0.040873	0.040873
{w1,w2}	0.4355082	0.338989	0.435508	0.096519
{w1,w3}	0.3980304	0.308165	0.398030	0.089866
{w1,w4}	0.2406251	0.187469	0.240625	0.053157
{w2,w3}	0.4546004	0.353964	0.454600	0.100637
{w2,w4}	0.2910446	0.233267	0.291045	0.057777
{w3,w4}	0.2571967	0.202443	0.257197	0.054754
{w1,w2,w3}	0.7545245	0.787580	0.787580	0.000000
{w1,w2,w4}	0.5538838	0.587315	0.587315	0.000000
{w1,w3,w4}	0.5133675	0.546814	0.546814	0.000000
{w2,w3,w4}	0.5745835	0.608005	0.608005	0.000000
{w1,w2,w3,w4}	1.0000000	0.994141	1.000000	0.005859

Table 6 — Discounted pseudo-belief function ($\delta = 0.02$)

Another way to restore feasibility is **Shafer’s discounting**.

Applying a discount factor $\delta = 0.02$ multiplies every proper subset by $(1 - \delta)$ while keeping $g(\Omega) = 1$. This uniform reduction weakens the dominance requirements across all levels of the lattice:

$$\hat{g}(A) = (1 - \delta)g(A), \quad A \neq \Omega.$$

The discounted function \hat{g} becomes compatible with a valid belief function, and the UAP again **succeeds**, reproducing Table 6 of the article.

These three examples highlight that the UAP is highly sensitive to the shape of the pseudo-belief function, and even minimal adjustments—either statistical (α) or structural (δ)—may determine whether a feasible upper approximation exists.

Table 6 - Discounting delta = 0.02 applied to Jeffreys g from Table 4

```
delta <- 0.02

# Start from g of Table 4 (alpha = 0.05)
g4_vals <- g4_res$g
names(g4_vals) <- g4_res$subset_names

# Apply Shafer discounting to all proper subsets
g6_vals <- g4_vals
is_top <- names(g6_vals) == "{w1,w2,w3,w4}"

g6_vals[!is_top] <- (1 - delta) * g6_vals[!is_top]
g6_vals[is_top] <- 1 # enforce g(Omega) = 1

# Run UAP
res6 <- upper_approximation(Omega4, g6_vals)

# Expected: TRUE
res6$success

## [1] TRUE

# Construct table
tab6 <- uap_table(res6, g6_vals)

# print table
knitr::kable(
  tab6,
  caption = "Table 6 - UAP succeeds after Shafer discounting with delta = 0.02.",
  align = "lrrrr",
  row.names = FALSE
)
```

Table 5: Table 6 — UAP succeeds after Shafer discounting with $\delta = 0.02$.

A	g	sum_prev	f	m_f
{w1}	0.1601868	0.000000	0.160187	0.160187
{w2}	0.2072201	0.000000	0.207220	0.207220
{w3}	0.1756192	0.000000	0.175619	0.175619

A	g	sum_prev	f	m_f
{w4}	0.0486323	0.000000	0.048632	0.048632
{w1,w2}	0.4513635	0.367407	0.451363	0.083957
{w1,w3}	0.4141842	0.335806	0.414184	0.078378
{w1,w4}	0.2562967	0.208819	0.256297	0.047478
{w2,w3}	0.4702490	0.382839	0.470249	0.087410
{w2,w4}	0.3072090	0.255852	0.307209	0.051357
{w3,w4}	0.2730705	0.224251	0.273071	0.048819
{w1,w2,w3}	0.7620079	0.792771	0.792771	0.000000
{w1,w2,w4}	0.5678782	0.598830	0.598830	0.000000
{w1,w3,w4}	0.5281538	0.559113	0.559113	0.000000
{w2,w3,w4}	0.5881103	0.619057	0.619057	0.000000
{w1,w2,w3,w4}	1.0000000	0.989056	1.000000	0.010944

Polyhedral Representation of the Upper Approximation

The Upper Approximation of pseudo-belief function g defined in the article is the set of all basic probability assignments m whose induced belief function f dominates the pseudo-belief function g .

For every subset $A \subseteq \Omega$,

$$f(A) = \sum_{B \subseteq A} m(B) \quad \text{and} \quad f(A) \geq g(A).$$

Together with the standard belief-function constraints

$$m(\emptyset) = 0, \quad m(A) \geq 0, \quad \sum_A m(A) = 1,$$

the feasible region becomes a **polytope** in $\mathbb{R}^{2^{|\Omega|}}$, described entirely by linear inequalities:

$$\mathcal{M}_g = \{m \mid \mathbf{M}_{\text{bel}} m \geq g, m \geq 0, m(\emptyset) = 0, \sum_A m(A) = 1\}.$$

Here, \mathbf{M}_{bel} is the *incidence matrix* of the subset lattice: its entry is 1 exactly when $B \subseteq A$. This matrix linearly maps a BPA m to the corresponding belief function f .

The polyhedral formulation serves two purposes:

1. it provides a global geometric interpretation of the dominance condition $f \geq g$;
2. it allows the use of dedicated tools (such as **cddlib**) to enumerate the **vertices** of \mathcal{M}_g and analyse its structure.

In particular, the ternary example in the article yields **28** extreme points, while the quaternary example expands to **13 889** vertices, illustrating the rapid combinatorial growth of the polytope as the dimension increases.

Constructing the Polytope in R

The following function constructs this polyhedron directly in terms of its linear inequality and equality constraints. This representation is *unified*: it is suitable both for **vertex enumeration** (**rcdd**) and for **linear programming** (LP-based approximations).

```
### Constructing the polytope  $M_g$  using a unified LP representation
build_lp_constraints <- function(Omega, g_values) {
```

```
  subsets      <- powerset(Omega)
  subset_names <- sapply(subsets, subset_name)
  n_vars       <- length(subsets)
```

```
  stopifnot(length(g_values) == n_vars)
```

```
  # -----
  # Dominance constraints:  $M_{bel} m \geq g$ 
  # -----
```

```
  M_dom <- matrix(0, nrow=n_vars, ncol=n_vars)
  for (i in seq_along(subsets)) {
    A <- subsets[[i]]
    M_dom[i, sapply(subsets, function(B) all(B %in% A))] <- 1
  }
  b_dom <- g_values
```

```
  # -----
  # Nonnegativity:  $m \geq 0$ 
  # -----
  M_nonneg <- diag(n_vars)
  b_nonneg <- rep(0, n_vars)
```

```
  # Stack all inequalities
  M_ineq <- rbind(M_dom, M_nonneg)
  b_ineq <- c(b_dom, b_nonneg)
```

```
  # -----
  # Equalities:
  #   sum m(A) = 1
  #   m(emptyset) = 0
  # -----
  M_eq <- matrix(0, nrow=2, ncol=n_vars)
  M_eq[1, ] <- 1           # normalization
  M_eq[2, 1] <- 1          # empty set mass = 0
  b_eq <- c(1, 0)
```

```
  colnames(M_ineq) <- subset_names
  colnames(M_eq)   <- subset_names
```

```
  list(
    subsets = subsets,
    subset_names = subset_names,
    M_ineq = M_ineq,
    b_ineq = b_ineq,
    M_eq = M_eq,
    b_eq = b_eq
  )
}
```

```
# Example build
```

```

Omega <- c("w1","w2","w3")
g_example <- c(0, 0.1,0.2,0.3, 0.5,0.6,0.8, 1.0)

lp_data <- build_lp_constraints(Omega, g_example)
lp_data[3:6]

## $M_ineq
##      {} {w1} {w2} {w3} {w1,w2} {w1,w3} {w2,w3} {w1,w2,w3}
## [1,] 1 0 0 0 0 0 0 0
## [2,] 1 1 0 0 0 0 0 0
## [3,] 1 0 1 0 0 0 0 0
## [4,] 1 0 0 1 0 0 0 0
## [5,] 1 1 1 0 1 0 0 0
## [6,] 1 1 0 1 0 1 0 0
## [7,] 1 0 1 1 0 0 1 0
## [8,] 1 1 1 1 1 1 1 1
## [9,] 1 0 0 0 0 0 0 0
## [10,] 0 1 0 0 0 0 0 0
## [11,] 0 0 1 0 0 0 0 0
## [12,] 0 0 0 1 0 0 0 0
## [13,] 0 0 0 0 1 0 0 0
## [14,] 0 0 0 0 0 1 0 0
## [15,] 0 0 0 0 0 0 1 0
## [16,] 0 0 0 0 0 0 0 1
##
## $b_ineq
## [1] 0.0 0.1 0.2 0.3 0.5 0.6 0.8 1.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0
##
## $M_eq
##      {} {w1} {w2} {w3} {w1,w2} {w1,w3} {w2,w3} {w1,w2,w3}
## [1,] 1 1 1 1 1 1 1 1
## [2,] 1 0 0 0 0 0 0 0
##
## $b_eq
## [1] 1 0

```

Vertex Enumeration of \mathcal{M}_g

The unified LP description of \mathcal{M}_g can be converted into the H -representation required by `rcdd`. Each vertex of this polytope corresponds to an *extreme* belief function that still dominates the pseudo-belief function g .

The number of vertices grows rapidly, as illustrated on polytopes related to above mentioned examples on $|\Omega| = 3, 4$. We also list first 10 vertices for each polytope.

This illustrates the exponential geometric complexity of the dominance constraints.

```

library(rcdd)

# Convert LP constraints to rcdd H-representation
lp_to_H <- function(lp) {
  H_ineq <- cbind(0, -lp$b_ineq, lp$M_ineq) # inequalities: 0 + Mx - b >= 0
  H_eq <- cbind(1, -lp$b_eq, lp$M_eq) # equalities: type=1
  d2q(rbind(H_ineq, H_eq))
}

```

```

# Extract vertices returned by rcdd
extract_vertices <- function(Vout) {
  out <- Vout$output
  V <- out[out[,1]=="0", , drop=FALSE] # rows marked as vertices
  if (nrow(V) == 0) return(NULL)
  q2d(V[, -c(1,2), drop=FALSE]) # drop type and constant
}

# Wrapper for computing vertices directly from g-values
enumerate_vertices <- function(Omega, g_values) {
  lp <- build_lp_constraints(Omega, g_values)
  H <- lp_to_H(lp)
  V <- scdd(H)
  verts <- extract_vertices(V)
  colnames(verts) <- names(g_values)
  list(
    subsets = lp$subset_names,
    vertices = verts,
    n_vertices = if (is.null(verts)) 0L else nrow(verts)
  )
}

# --- Ternary example ---
res <- enumerate_vertices(Omega3, g_ternary)
res$n_vertices # expected: 28

```

```
## [1] 28
```

```
res$vertices[1:10, ]
```

```

##      {} {w1} {w2} {w3}      {w1,w2} {w1,w3} {w2,w3} {w1,w2,w3}
## [1,] 0  0.1  0.4  0.3 5.551115e-17      0.0      0.0      0.2
## [2,] 0  0.1  0.2  0.5 2.000000e-01      0.0      0.0      0.0
## [3,] 0  0.2  0.5  0.2 1.000000e-01      0.0      0.0      0.0
## [4,] 0  0.1  0.4  0.3 2.000000e-01      0.0      0.0      0.0
## [5,] 0  0.2  0.2  0.2 1.000000e-01      0.0      0.3      0.0
## [6,] 0  0.1  0.2  0.3 2.000000e-01      0.0      0.2      0.0
## [7,] 0  0.1  0.3  0.2 1.000000e-01      0.1      0.2      0.0
## [8,] 0  0.1  0.5  0.2 1.000000e-01      0.1      0.0      0.0
## [9,] 0  0.1  0.4  0.2 0.000000e+00      0.1      0.1      0.1
## [10,] 0  0.1  0.4  0.2 0.000000e+00      0.2      0.1      0.0

```

```

# --- Quaternary example ---
res4 <- enumerate_vertices(Omega4, g5_vals)
res4$n_vertices # expected: 13889

```

```
## [1] 13889
```

```
res4$vertices[1:10, ]
```

```

##      {}      {w1}      {w2}      {w3}      {w4}      {w1,w2} {w1,w3} {w1,w4}
## [1,] 0 0.1997516 0.2501711 0.2415141 0.04087347 0.06308764      0      0
## [2,] 0 0.1997516 0.1923940 0.2992912 0.04087347 0.06308764      0      0
## [3,] 0 0.1465951 0.1923940 0.3190163 0.11837567 0.09651909      0      0
## [4,] 0 0.1465951 0.1923940 0.3190163 0.24547552 0.09651909      0      0
## [5,] 0 0.1465951 0.1923940 0.3190163 0.11837567 0.09651909      0      0

```



```

## [6,] 0 0.1465951 0.1923940 0.3190163 0.09402999 0.09651909 0 0
## [7,] 0 0.1465951 0.1923940 0.3190163 0.09402999 0.09651909 0 0
## [8,] 0 0.1465951 0.1923940 0.2727424 0.09402999 0.14279295 0 0
## [9,] 0 0.1465951 0.1923940 0.2727424 0.09402999 0.09651909 0 0
## [10,] 0 0.1465951 0.1923940 0.2727424 0.09402999 0.09651909 0 0
##      {w2,w3}      {w2,w4}      {w3,w4} {w1,w2,w3} {w1,w2,w4} {w1,w3,w4} {w2,w3,w4}
## [1,]      0 0.00000000 0.0000000 0.00000000      0 0.03122834 0.1733737
## [2,]      0 0.05777713 0.0000000 0.00000000      0 0.14682492 0.0000000
## [3,]      0 0.00000000 0.0000000 0.00000000      0 0.00000000 0.1270999
## [4,]      0 0.00000000 0.0000000 0.00000000      0 0.00000000 0.0000000
## [5,]      0 0.00000000 0.1270999 0.00000000      0 0.00000000 0.0000000
## [6,]      0 0.02434568 0.0000000 0.00000000      0 0.00000000 0.1270999
## [7,]      0 0.15144554 0.0000000 0.00000000      0 0.00000000 0.0000000
## [8,]      0 0.15144554 0.0000000 0.00000000      0 0.00000000 0.0000000
## [9,]      0 0.02434568 0.0000000 0.04627386      0 0.00000000 0.1270999
## [10,]      0 0.15144554 0.0000000 0.04627386      0 0.00000000 0.0000000
##      {w1,w2,w3,w4}
## [1,]      0
## [2,]      0
## [3,]      0
## [4,]      0
## [5,]      0
## [6,]      0
## [7,]      0
## [8,]      0
## [9,]      0
## [10,]      0

```

Linear Programming Approximations

The polytope \mathcal{M}_g contains all belief functions f dominating the pseudo-belief function g . To select a *single representative* belief function, we minimise a linear objective

$$\min_{m \in \mathcal{M}_g} c^\top m,$$

where the vector c encodes a preference for certain focal sets.

We now implement three criteria used in the article:

- **L1** (minimal upward correction)
- **HD** (negative Dubois–Prade entropy; LP minimisation = entropy maximisation)
- **CW** (cardinality-weighted)

The following unified solver builds the LP constraints and computes the optimal m and induced belief function f for any objective.

```

objective_L1 <- function(subsets, Omega) {
  sapply(subsets, function(B) 2^(length(Omega) - length(B)))
}

objective_HD <- function(subsets, Omega) {
  - sapply(subsets, function(B) if (length(B) == 0) 0 else log(length(B), base = 2))
}

```

```
objective_CW <- function(subsets, Omega) {
  sapply(subsets, function(B) if (length(B) == 0) 0 else 1/length(B))
}
```

Call the LP solver

```
# =====
# solve_lp_for_objective()
# -----
# Solves the linear program defining the upper approximation  $f \geq g$ 
# for a chosen linear objective function in the belief space.
#
# The LP is formulated over the basic probability assignment vector  $m$ ,
# subject to the standard belief-function constraints:
#
# (1)  $M_{bel} * m \geq g$  (dominance:  $f(A) \geq g(A)$ )
# (2)  $m(A) \geq 0$  (non-negativity)
# (3)  $m(\text{emptyset}) = 0$  (empty set has zero mass)
# (4)  $\sum_A m(A) = 1$  (normalization)
#
# The objective is always of the form:
# minimize  $\sum_A w(A) * m(A)$ 
# where  $w(A)$  is supplied by obj_fun (L1, HD, CW, ...).
#
# Input:
# Omega ... frame of discernment
# g_values ... pseudo-belief function  $g(A)$ 
# obj_fun ... objective-generating function, e.g. objective_L1
#
# Output:
# A list with:
# subset ... subset names in lexicographic order
# f ... resulting belief function  $f(A)$ 
# m ... resulting BPA  $m(A)$ 
#
# =====
solve_lp_for_objective <- function(Omega, g_values, obj_fun) {

  # Construct all LP constraint matrices for  $f \geq g$ 
  lp_data <- build_lp_constraints(Omega, g_values)

  subsets <- lp_data$subsets
  subset_names <- lp_data$subset_names
  nA <- length(subsets)

  # -----
  # Objective vector:  $w(A)$  for each subset  $A$ 
  # (determined by objective_L1, objective_HD, objective_CW, ...)
  # -----
  obj <- obj_fun(subsets, Omega)

  # -----
  # Full LP system:
```

```

# [M_ineq] m >= b_ineq
# [M_eq ] m == b_eq
# -----
M_full <- rbind(lp_data$M_ineq, lp_data$M_eq)
dir_full <- c(rep(">=", nrow(lp_data$M_ineq)), rep("=", nrow(lp_data$M_eq)))
rhs_full <- c(lp_data$b_ineq, lp_data$b_eq)

# -----
# Solve LP using lpSolve
# -----
sol <- lp(
  direction = "min",
  objective.in = obj,
  const.mat = M_full,
  const.dir = dir_full,
  const.rhs = rhs_full
)

# Extract BPA m(A)
m <- sol$solution
names(m) <- subset_names

# Reconstruct belief f(A) = sum_{B \subseteq A} m(B)
f <- sapply(subsets, function(A) {
  idx <- sapply(subsets, function(B) all(B %in% A))
  sum(m[idx])
})

list(
  subset = subset_names,
  f = f,
  m = m
)
}

```

LP Reproduction of Tables 7, 8, 9

The following code block reproduces Tables 7, 8, and 9 from the paper by:

1. constructing the polytope \mathcal{M}_g ,
2. solving the LP for each objective,
3. computing the corresponding belief and mass assignments,
4. presenting all results in a unified table.

As in the article, we omit the empty set from printed tables.

```

# =====
# compute_table()
# -----
# Computes all mass and belief values for several approximation
# methods (L1, CW, HD, UAP) and returns a unified table.
#
# Input:
# Omega      - vector of states ("w1", "w2", ...)
# g_values   - pseudo-belief function g(A), already ordered and normalized

```

```

#
# Output:
#   A list with:
#     $table - data frame containing g(A), f(A), and m(A) for all methods
#     $subsets - list of subsets of Omega (used by criteria function)
#
# This function does NOT compute criterion values; that is delegated
# to compute_criteria(), keeping responsibilities cleanly separated.
# =====
compute_table <- function(Omega, g_values) {

  lp_data <- build_lp_constraints(Omega, g_values)
  subsets <- lp_data$subsets
  subset_names <- lp_data$subset_names
  idx <- 2:length(subset_names) # remove empty set

  # ---- Solve all LP variants ----
  sol_L1 <- solve_lp_for_objective(Omega, g_values, objective_L1)
  sol_HD <- solve_lp_for_objective(Omega, g_values, objective_HD)
  sol_CW <- solve_lp_for_objective(Omega, g_values, objective_CW)
  uap <- upper_approximation(Omega, g_values)

  # ---- Build main table (body) ----
  # Each row corresponds to a subset A in lexicographic order.
  # Columns contain:
  #   g(A) - pseudo-belief
  #   f_method - belief produced by each approximation method
  #   m_method - corresponding mass assignments
  table_body <- data.frame(
    subset = subset_names,
    g = g_values,

    f_L1 = sol_L1$f,
    f_HD = sol_HD$f,
    f_CW = sol_CW$f,
    f_UAP = uap$f,

    m_L1 = sol_L1$m,
    m_HD = sol_HD$m,
    m_CW = sol_CW$m,
    m_UAP = uap$m,

    row.names = NULL
  )

  list(
    table_body = table_body,
    subsets = subsets
  )
}

# =====
# evaluate_lp_criteria()

```

```

# -----
# Computes values of each criterion (L1, CW, HD) for each mass
# vector in the table produced by compute_table_explicit().
#
# Input:
#   table_body - data frame containing m_* columns
#   subsets    - list of subsets (same order as masses)
#   Omega      - frame of discernment
#
# Output:
#   A matrix with rows = criteria (L1, CW, HD)
#                   columns = mass vectors (m_L1, m_CW, m_HD, m_rSP, m_UAP ...)
#
# The formula is always:
#   criterion_value = sum_A w(A) * m(A)
# where w(A) is given by objective_* functions.
# =====
evaluate_lp_criteria <- function(table_body, subsets, Omega) {

  # Find all mass columns automatically (m_L1, m_CW, m_HD, m_rSP, m_UAP, ...)
  m_cols <- grep("^m_", names(table_body), value = TRUE)

  # Available criteria (objective functions defined earlier)
  crit_map <- list(
    L1 = objective_L1,
    CW = objective_CW,
    HD = function(subsets, Omega) {-objective_HD(subsets, Omega)}
  )

  # Initialize (criteria × mass vectors)
  crit_matrix <- matrix(
    NA,
    nrow = length(crit_map),
    ncol = length(m_cols),
    dimnames = list(names(crit_map), m_cols)
  )

  # Calculate weighted sum for each criterion-method pair
  for (crit_name in names(crit_map)) {
    w <- crit_map[[crit_name]](subsets, Omega) # objective weights
    for (m_col in m_cols) {
      m_vec <- table_body[[m_col]]
      crit_matrix[crit_name, m_col] <- sum(w * m_vec)
    }
  }

  crit_matrix
}

```

Table 7

```

table <- compute_table(g_values = g_ternary, Omega = Omega3)
knitr::kable(

```

```

table$table_body[-1,],
digits = 3,
caption = "Upper approximations (L1, CW, HD, UAP) for the ternary example  $g$ "
)

```

Table 6: Upper approximations (L1, CW, HD, UAP) for the ternary example g

	subset	g	f_L1	f_HD	f_CW	f_UAP	m_L1	m_HD	m_CW	m_UAP
2	{w1}	0.1	0.2	0.1	0.2	0.1	0.2	0.1	0.2	0.1
3	{w2}	0.1	0.2	0.3	0.2	0.1	0.2	0.3	0.2	0.1
4	{w3}	0.2	0.2	0.2	0.2	0.2	0.2	0.2	0.2	0.2
5	{w1,w2}	0.5	0.5	0.5	0.5	0.5	0.1	0.1	0.1	0.3
6	{w1,w3}	0.4	0.4	0.4	0.4	0.4	0.0	0.1	0.0	0.1
7	{w2,w3}	0.7	0.7	0.7	0.7	0.7	0.3	0.2	0.3	0.4
8	{w1,w2,w3}	1.0	1.0	1.0	1.0	1.2	0.0	0.0	0.0	0.0

```

evaluate_lp_criteria(table_body = table$table_body, subsets = table$subsets, Omega = Omega3)

```

```

##      m_L1 m_HD m_CW m_UAP
## L1  3.2  3.2  3.2  3.2
## CW  0.8  0.8  0.8  0.8
## HD  0.4  0.4  0.4  0.8

```

Table 8

```

table <- compute_table(g_values = g_ternary_tilde, Omega = Omega3)
knitr::kable(
  table$table_body[-1,],
  digits = 3,
  caption = "Upper approximations (L1, CW, HD, UAP) for the ternary example  $\tilde{g}$ "
)

```

Table 7: Upper approximations (L1, CW, HD, UAP) for the ternary example \tilde{g}

	subset	g	f_L1	f_HD	f_CW	f_UAP	m_L1	m_HD	m_CW	m_UAP
2	{w1}	0.2	0.2	0.2	0.2	0.2	0.2	0.2	0.2	0.2
3	{w2}	0.2	0.2	0.2	0.2	0.2	0.2	0.2	0.2	0.2
4	{w3}	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3	0.3
5	{w1,w2}	0.5	0.5	0.5	0.5	0.5	0.1	0.1	0.1	0.1
6	{w1,w3}	0.4	0.5	0.5	0.5	0.5	0.0	0.0	0.0	0.0
7	{w2,w3}	0.7	0.7	0.7	0.7	0.7	0.2	0.2	0.2	0.2
8	{w1,w2,w3}	1.0	1.0	1.0	1.0	1.0	0.0	0.0	0.0	0.0

```

evaluate_lp_criteria(table_body = table$table_body, subsets = table$subsets, Omega = Omega3)

```

```

##      m_L1 m_HD m_CW m_UAP
## L1  3.40  3.40  3.40  3.40
## CW  0.85  0.85  0.85  0.85
## HD  0.30  0.30  0.30  0.30

```

Table 9

```
table <- compute_table(g_values = g5_vals, Omega = Omega4)
knitr::kable(
  table$table_body[-1,],
  digits = 3,
  caption = "Upper approximations (L1, CW, HD, UAP) for the quaternary example"
)
```

Table 8: Upper approximations (L1, CW, HD, UAP) for the quaternary example

	subset	g	f_L1	f_HD	f_CW	f_UAP	m_L1	m_HD	m_CW	m_UAP
2	{w1}	0.147	0.158	0.158	0.147	0.147	0.158	0.158	0.147	0.147
3	{w2}	0.192	0.203	0.203	0.192	0.192	0.203	0.203	0.192	0.192
4	{w3}	0.162	0.173	0.173	0.162	0.162	0.173	0.173	0.162	0.162
5	{w4}	0.041	0.052	0.052	0.041	0.041	0.052	0.052	0.041	0.041
6	{w1,w2}	0.436	0.436	0.436	0.436	0.436	0.074	0.074	0.097	0.097
7	{w1,w3}	0.398	0.398	0.398	0.398	0.398	0.068	0.068	0.090	0.090
8	{w1,w4}	0.241	0.241	0.241	0.241	0.241	0.031	0.031	0.053	0.053
9	{w2,w3}	0.455	0.455	0.455	0.455	0.455	0.079	0.079	0.101	0.101
10	{w2,w4}	0.291	0.291	0.291	0.291	0.291	0.035	0.035	0.058	0.058
11	{w3,w4}	0.257	0.257	0.257	0.257	0.257	0.032	0.032	0.055	0.055
12	{w1,w2,w3}	0.755	0.755	0.755	0.788	0.788	0.000	0.000	0.000	0.000
13	{w1,w2,w4}	0.554	0.554	0.554	0.587	0.587	0.000	0.000	0.000	0.000
14	{w1,w3,w4}	0.513	0.513	0.513	0.547	0.547	0.000	0.000	0.000	0.000
15	{w2,w3,w4}	0.575	0.575	0.575	0.608	0.608	0.000	0.000	0.000	0.000
16	{w1,w2,w3,w4}	1.000	1.000	1.000	1.000	1.000	0.095	0.095	0.006	0.006

```
evaluate_lp_criteria(table_body = table$table_body, subsets = table$subsets, Omega = Omega4)
```

```
##           m_L1      m_HD      m_CW      m_UAP
## L1 6.0592484 6.0592484 6.1481516 6.1481516
## CW 0.7692513 0.7692513 0.7692513 0.7692513
## HD 0.5088785 0.5088785 0.4644269 0.4644269
```

Table 10

Number of vertices with the same value

```
# =====
# compute_vertex_minima()
# -----
# Computes, for a pseudo-belief function g, how many extreme points
# of the polytope M_g minimize each LP criterion (L1, HD, CW).
#
# The function:
# 1. constructs the polytope M_g via enumerate_vertices(),
# 2. enumerates all extreme points (vertices),
# 3. evaluates each LP objective over all vertices,
# 4. counts how many vertices achieve the minimum value.
#
# Input:
# g_values ... numeric vector of g(A) in the same subset order
```

```

# Omega    ... frame of discernment, e.g. c("w1","w2","w3")
#
# Output:
# data.frame with one row and columns:
#   all_vertices ... total number of extreme points
#   L1           ... count of vertices minimizing L1 objective
#   HD           ... count of vertices minimizing HD objective
#   CW           ... count of vertices minimizing CW objective
#
# The result corresponds to Table 10 in the article.
# =====
compute_vertex_minima <- function(g_values, Omega) {

  # enumerate all vertices of M_g
  verts <- enumerate_vertices(Omega, g_values)$vertices
  subsets <- powerset(Omega)

  # number of vertices
  n_verts <- nrow(verts)

  # criterion → objective function
  crit_map <- list(
    L1 = objective_L1,
    HD = objective_HD,
    CW = objective_CW
  )

  # initialize result vector
  out <- c(
    all_vertices = n_verts,
    L1 = NA,
    HD = NA,
    CW = NA
  )

  # compute for each criterion
  for (crit in names(crit_map)) {
    w <- crit_map[[crit]](subsets, Omega)
    vals <- verts %*% w
    min_val <- min(vals)
    out[crit] <- sum(abs(vals - min_val) < 1e-14)
  }

  as.data.frame(t(out))
}

## =====
## Construct Table 10 - counts of vertices minimizing each criterion
## =====

row1 <- compute_vertex_minima(g_ternary,      Omega3)
row2 <- compute_vertex_minima(g_ternary_tilde, Omega3)

```



```

row3 <- compute_vertex_minima(g5_vals, Omega4)

# Add descriptive labels (first column)
Table10 <- rbind(
  "Ternary original g"= row1,
  "Ternary modified g_tilde" = row2,
  "Quaternary example"= row3
)

# Convert row names into a proper column
Table10 <- cbind(
  Example = rownames(Table10),
  Table10,
  row.names = NULL
)

# Print table
knitr::kable(
  Table10,
  caption = "Table 10 - Counts of extreme points and numbers achieving minimum value under each objective",
  align = "lcr",
  digits = 0
)

```

Table 9: Table 10 — Counts of extreme points and numbers achieving minimum value under each objective.

Example	all_vertices	L1	HD	CW
Ternary original g	28	3	3	3
Ternary modified g_tilde	27	4	4	4
Quaternary example	13889	1	1	14