

4. Build a simulation manager to run each simulation.
5. Build a results manager to generate summary results (metrics).
6. Build a validator to select a model ensemble.

Step 1: Build the population model for the study region

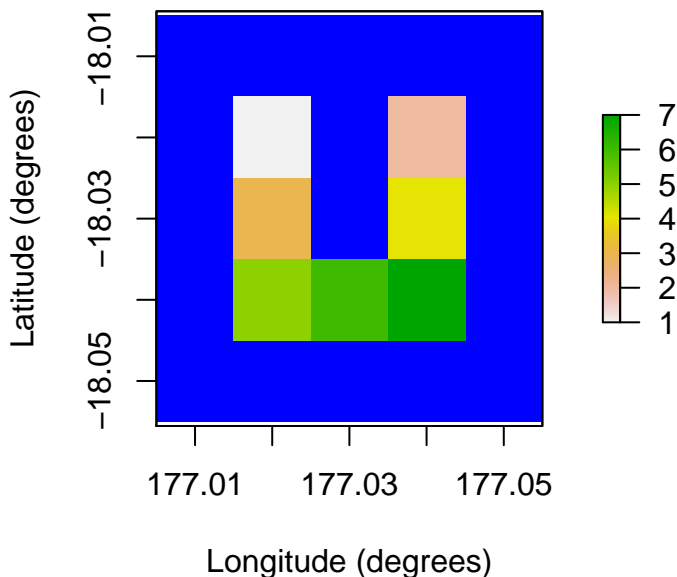
Create a model template using the *PopulationModel* class. If the model is spatially explicit, then define the study region via the *Region* class. All fixed model inputs, such as environmental correlations, as well as any user-defined functions for processes such as harvesting, should be set at this stage.

We could create these components in any order and set model parameters separately, but let's setup our study region, generate environmental correlations, and define a harvest function prior to initializing the template model with all the fixed parameters.

Study region First, we'll define our study region (denoted U Island) with some longitude-latitude coordinates.

```
# Demonstration example region (U Island)
coordinates <- data.frame(x = rep(seq(177.01, 177.05, 0.01), 5),
                          y = rep(seq(-18.01, -18.05, -0.01), each = 5))
template_raster <- Region$new(coordinates = coordinates)$region_raster # full extent
template_raster[[-c(7, 9, 12, 14, 17:19)] <- NA # make U Island
region <- Region$new(template_raster = template_raster)
raster::plot(region$region_raster, main = "Example region (cell indices)",
             xlab = "Longitude (degrees)", ylab = "Latitude (degrees)",
             colNA = "blue")
```

Example region (cell indices)



Environmental correlation Next, we'll define a distance-based spatial correlation for applying environmental stochasticity within our model. The generated correlation data is compacted for computational efficiency (with large-scale models).

```
# Distance-based environmental correlation (via a compacted Cholesky decomposition)
env_corr <- SpatialCorrelation$new(region = region, amplitude = 0.4, breadth = 500)
```

```

correlation <- env_corr$get_compact_decomposition(decimals = 2)
correlation # examine
#> $matrix
#>      [,1] [,2] [,3] [,4] [,5] [,6] [,7]
#> [1,]  1 0.01 0.04 0.04 0.04 0.02 0.04
#> [2,]  0 1.00 1.00 0.01 1.00 0.02 0.01
#> [3,]  0 0.00 0.00 1.00 0.00 0.05 0.05
#> [4,]  0 0.00 0.00 0.00 0.00 1.00 1.00
#>
#> $map
#>      [,1] [,2] [,3] [,4] [,5] [,6] [,7]
#> [1,]  1   1   1   2   3   3   4
#> [2,] NA   2   3   3   5   4   5
#> [3,] NA  NA  NA   4  NA   5   6
#> [4,] NA  NA  NA  NA  NA   6   7

```

Harvest function Let's now define a simple harvest function, which is optionally list-nested with a harvest rate parameter. We'll also define an alias to the harvest rate so we can sample this parameter later.

```

# User-defined harvest function (list-nested) and alias
harvest <- list(rate = NA, # sample later
               function(params) round(params$stage_abundance*(1 - params$rate)))
harvest_rate_alias <- list(harvest_rate = "harvest$rate")

```

Template model Finally, we can build our template model with these and other fixed parameters.

```

# Population (simulation) model template for fixed parameters
stage_matrix <- matrix(c(0, 2.5, # Leslie/Lefkovitch matrix
                       0.8, 0.5), nrow = 2, ncol = 2, byrow = TRUE,
                      dimnames = list(c("juv", "adult"), c("juv", "adult")))
stage_matrix # examine
#>      juv adult
#> juv  0.0  2.5
#> adult 0.8  0.5
model_template <- PopulationModel$new(region = region,
                                     time_steps = 10, # years
                                     populations = region$region_cells, # 7
                                     stages = 2,
                                     stage_matrix = stage_matrix,
                                     demographic_stochasticity = TRUE,
                                     standard_deviation = 0.05,
                                     correlation = correlation,
                                     density_dependence = "logistic",
                                     harvest = harvest,
                                     results_selection = c("abundance", "harvested"),
                                     attribute_aliases = harvest_rate_alias)

```

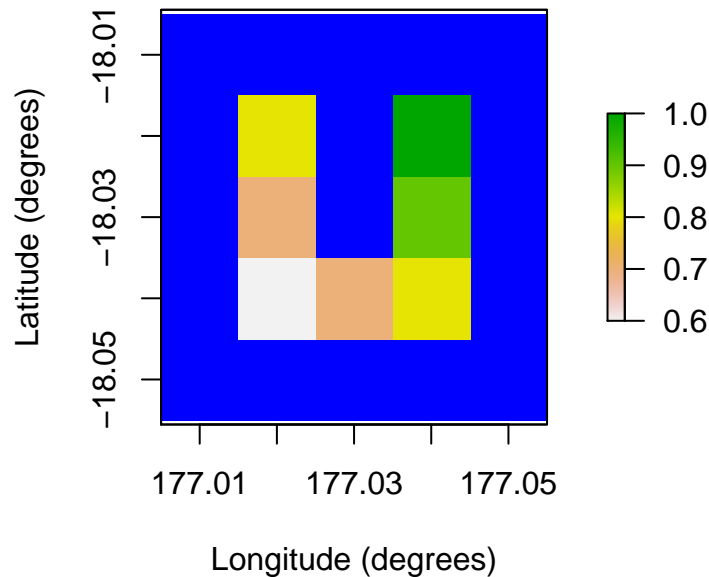
Step 2: Build generators for dynamically generating model parameters

Some model parameters are single values, whilst others are represented as arrays and other multi-value data structures. Usually, we don't wish to sample each individual value within these multi-value parameters (in step 3), but to generate them dynamically via one or more intermediate sampled parameters. Here we build generators for model initial abundance, carrying capacity, and dispersal.

Habitat suitability Firstly, our initial abundance and carrying capacity generator utilizes an example (mock) habitat suitability for our defined study region.

```
# Example habitat suitability
example_hs <- c(0.8, 1, 0.7, 0.9, 0.6, 0.7, 0.8 )
example_hs_raster <- region$region_raster
example_hs_raster[region$region_indices] <- example_hs
raster::plot(example_hs_raster, main = "Example habitat suitability",
             xlab = "Longitude (degrees)", ylab = "Latitude (degrees)",
             colNA = "blue")
```

Example habitat suitability



Initial abundance and carrying capacity generator The generator utilizes generic template functionality for user-defined custom functions. It uses sampled input parameters (initial total abundance and maximum cell density), along with habitat suitability, to generate the desired output model parameters (initial abundance and carrying capacity) via these user-defined functions. Generators can also be configured to read values from files or generate values via probabilistic distributions.

```
# Initial abundance and carrying capacity generated via example habitat suitability
capacity_gen <- Generator$new(description = "Capacity generator",
                             example_hs = example_hs, # template attached
                             inputs = c("initial_n", "density_max"),
                             outputs = c("initial_abundance", "carrying_capacity"))
capacity_gen$add_generative_requirements(list(initial_abundance = "function",
                                              carrying_capacity = "function"))
capacity_gen$add_function_template("initial_abundance",
                                   function_def = function(params) {
                                     stats::rmultinom(1, size = params$initial_n,
                                                       prob = params$example_hs)[,1]
                                   },
                                   call_params = c("initial_n", "example_hs"))
capacity_gen$add_function_template("carrying_capacity",
                                   function_def = function(params) {
                                     round(params$density_max*params$example_hs)
                                   },
```

```

                                call_params = c("density_max", "example_hs"))
capacity_gen$generate(input_values = list(initial_n = 500, density_max = 100)) # test
#> $initial_abundance
#> [1] 76 85 68 78 64 62 67
#>
#> $carrying_capacity
#> [1] 80 100 70 90 60 70 80

```

Dispersal generator Our dispersal generator uses default functionality for generating dispersal rates between cells. Its sampled inputs parameterize the distance-based dispersal function (proportion dispersing and breadth of dispersal). The generator can be configured with a dispersal friction helper class object, which calculates equivalent dispersal distances for frictional landscapes and coastlines. Here it is used to ensure dispersal is not performed directly across the “water” in our example U-shaped island. The generated dispersal data is calculated via pre-calculated distance data and compacted for computational efficiency (with large-scale models).

```

# Distance-based dispersal generator
dispersal_gen <- DispersalGenerator$new(region = region,
                                       dispersal_max_distance = 3000, # in m
                                       dispersal_friction = DispersalFriction$new(),
                                       inputs = c("dispersal_p", "dispersal_b"),
                                       decimals = 5)
dispersal_gen$calculate_distance_data() # pre-calculate
test_dispersal <- dispersal_gen$generate(input_values = list(dispersal_p = 0.5,
                                                            dispersal_b = 700))
head(test_dispersal$dispersal_data[[1]])
#>   target_pop source_pop emigrant_row immigrant_row dispersal_rate
#> 1           3           1             1             1         0.10284
#> 2           5           1             2             1         0.02115
#> 3           6           1             3             1         0.01501
#> 4           4           2             1             1         0.10284
#> 5           6           2             2             2         0.01501
#> 6           7           2             3             1         0.02115

```

Note that there is no dispersal rate between cells 1 and 2 as there is “water” between those cells, and consequently dispersal between those cells must travel around the U-shaped island, which can’t be achieved in one simulation time-step since the “round” distance between those cells is greater than 3000 m.

Step 3: Sample model and generator parameters for each simulation

In order to explore the model parameter space to find the best models, we generate Latin hypercube samples of model and generator parameters to be simulated, using the *LatinHypercubeSampler* class. This class has functionality for generating sample parameters via Uniform, Normal, Lognormal, Beta, and Triangular distributions. For our example we only generate 12 samples. We encourage the user to generate hundreds, or thousands, of samples.

```

# Generate sampled values for variable model parameters via LHS
lhs_gen <- LatinHypercubeSampler$new()
lhs_gen$set_uniform_parameter("growth_rate_max", lower = 0.4, upper = 0.6, decimals = 2)
lhs_gen$set_uniform_parameter("harvest_rate", lower = 0.05, upper = 0.15, decimals = 2)
lhs_gen$set_uniform_parameter("initial_n", lower = 400, upper = 600, decimals = 0)
lhs_gen$set_uniform_parameter("density_max", lower = 80, upper = 120, decimals = 0)
lhs_gen$set_uniform_parameter("dispersal_p", lower = 0.2, upper = 0.5, decimals = 2)
lhs_gen$set_uniform_parameter("dispersal_b", lower = 400, upper = 1000, decimals = 0)
sample_data <- lhs_gen$generate_samples(number = 12, random_seed = 123)
sample_data # examine

```

```

#>   growth_rate_max harvest_rate initial_n density_max dispersal_p dispersal_b
#> 1      0.50      0.09      575      111      0.26      481
#> 2      0.41      0.07      487      96      0.34      589
#> 3      0.44      0.09      483      110      0.42      509
#> 4      0.59      0.11      589      86      0.38      859
#> 5      0.56      0.06      441      92      0.21      667
#> 6      0.51      0.14      458      116      0.50      980
#> 7      0.54      0.05      416      88      0.28      447
#> 8      0.43      0.13      542      107      0.36      920
#> 9      0.46      0.12      422      101      0.43      849
#> 10     0.52      0.08      552      99      0.32      745
#> 11     0.48      0.11      525      119      0.25      789
#> 12     0.58      0.15      505      81      0.45      609

```

Step 4: Build a simulation manager to run each simulation

We now wish to run a simulation for each set (or row) of sampled parameters. The *SimulationManager* class manages the generation of parameters (via the generators), the running the model simulations, and writing simulation results to disk. It also maintains a log of each simulation's success and any errors or warnings encountered.

```

# Create a simulation manager and run the sampled model simulations
sim_manager <- SimulationManager$new(sample_data = sample_data,
                                     model_template = model_template,
                                     generators = list(capacity_gen, dispersal_gen),
                                     parallel_cores = 2,
                                     results_dir = OUTPUT_DIR)

run_output <- sim_manager$run()
run_output$summary
#> [1] "12 of 12 sample models ran and saved results successfully"
dir(OUTPUT_DIR, "*.RData") # includes 12 result files
#> [1] "sample_10_results.RData" "sample_11_results.RData"
#> [3] "sample_12_results.RData" "sample_1_results.RData"
#> [5] "sample_2_results.RData" "sample_3_results.RData"
#> [7] "sample_4_results.RData" "sample_5_results.RData"
#> [9] "sample_6_results.RData" "sample_7_results.RData"
#> [11] "sample_8_results.RData" "sample_9_results.RData"
dir(OUTPUT_DIR, "*.txt") # plus simulation log
#> [1] "simulation_log.txt"

```

Note that the output directory contains a R-data result files for each sample simulation and a simulation log file.

Step 5: Build a results manager to generate summary results (metrics)

We now wish to collate summary results for each of our simulations via the *ResultsManager* class. This manager loads the results from each sample simulation into an intermediate *PopulationResults* class object, which dynamically generates further results. We need to define functions for calculating summary metrics, as well as any matrices (one row of values per simulation) that we may be interested in examining. Each metric (or matrix) is associated with a user-defined function that utilizes results object attributes, or alternatively direct access to an attribute may be defined via a string. Once generated, the result metrics (a data frame) and/or matrices (a list) can be accessed via the manager. We may utilize the collated results in a variety of ways. However, with the objective of selecting the best models, we wish to compare (or validate) these result metrics to (with) known or desired target patterns (in step 6).

```

results_manager <- ResultsManager$new(simulation_manager = sim_manager,
                                     simulation_results = PopulationResults$new(),
                                     summary_metrics = c("trend_n", "total_h"),
                                     summary_matrices = c("n", "h"),
                                     summary_functions = list(
                                       trend_n = function(results) {
                                         round(results$all$abundance_trend, 2)
                                       },
                                       total_h = function(results) {
                                         sum(results$harvested)
                                       },
                                       n = "all$abundance", # string
                                       h = "all$harvested"),
                                     parallel_cores = 2)

gen_output <- results_manager$generate()
gen_output$summary
#> [1] "12 of 12 summary metrics/matrices generated from sample results successfully"
dir(OUTPUT_DIR, "*.txt") # plus generation log
#> [1] "generation_log.txt" "simulation_log.txt"
results_manager$summary_metric_data
#>   index trend_n total_h
#> 1     1     1.25    495
#> 2     2    -9.75    328
#> 3     3    -1.33    463
#> 4     4    -0.29    481
#> 5     5     4.60    294
#> 6     6     0.60    735
#> 7     7     2.75    226
#> 8     8   -10.71    635
#> 9     9     2.75    521
#> 10    10    -3.83    403
#> 11    11    -3.00    630
#> 12    12    -0.33    589
results_manager$summary_matrix_list
#> $n
#>   [,1] [,2] [,3] [,4] [,5] [,6] [,7] [,8] [,9] [,10]
#> [1,] 519 502 464 483 471 494 468 488 535 512
#> [2,] 457 488 451 463 424 432 441 410 377 393
#> [3,] 468 483 452 457 498 467 483 447 470 449
#> [4,] 457 381 373 358 357 355 368 400 379 425
#> [5,] 458 439 459 468 479 449 495 494 491 445
#> [6,] 457 434 486 425 449 431 467 438 449 452
#> [7,] 422 442 428 438 441 442 463 436 421 464
#> [8,] 477 488 417 431 425 420 411 402 391 342
#> [9,] 400 415 353 357 367 379 381 374 378 398
#> [10,] 490 459 473 452 473 430 427 436 448 512
#> [11,] 520 539 493 511 498 490 472 526 490 503
#> [12,] 390 344 333 311 321 353 316 349 331 354
#>
#> $h
#>   [,1] [,2] [,3] [,4] [,5] [,6] [,7] [,8] [,9] [,10]
#> [1,] 52 49 48 49 47 49 47 49 52 53
#> [2,] 35 37 35 34 31 33 33 32 28 30
#> [3,] 46 47 45 47 51 47 48 40 46 46
#> [4,] 56 49 45 46 45 45 46 51 46 52
#> [5,] 28 27 29 29 30 27 32 31 32 29

```

```

#> [6,] 76 71 79 69 76 71 76 71 73 73
#> [7,] 23 20 23 23 23 24 23 22 21 24
#> [8,] 71 73 62 67 63 64 63 61 61 50
#> [9,] 56 56 48 48 50 52 52 53 52 54
#> [10,] 42 41 43 38 42 37 37 39 38 46
#> [11,] 65 65 60 66 61 63 60 65 62 63
#> [12,] 67 61 57 55 56 61 53 60 57 62

```

Step 6: Build a validator to select a model ensemble

We now select and analyze our ‘best’ models via a *Validator* class object, which by default utilizes an approximate Bayesian computation (ABC) approach (Beaumont, Zhang, & Balding, 2002) provided by the *abc* library (Csillery et al., 2015). The validator can be configured appropriately for a problem domain (see the *abc* documentation). Here we use the default configuration to select the best 3 models, along with a weight value, which is indicative of the congruence between each model’s summary metrics and the corresponding target patterns. Also provided (with the default settings) is a diagnostic output (PDF) file, containing statistical information for analyzing the contribution of model parameters in the selection/validation process (see the *abc* documentation). For our simple demonstration the metrics and corresponding targets are relatively trivial, having the aim of producing stable population abundances with high harvest. However, the package facilitates the use more complex spatio-temporal metrics and targets (demonstrated in more advanced vignettes).

```

# Create a validator for selecting the 'best' example models
validator <- Validator$new(simulation_parameters = sample_data,
                          simulation_summary_metrics =
                            results_manager$summary_metric_data[-1],
                          observed_metric_targets = c(trend_n = 0, total_h = 600),
                          output_dir = OUTPUT_DIR)
suppressWarnings(validator$run(tolerance = 0.25, output_diagnostics = TRUE))
#> 12345678910
#> 12345678910
dir(OUTPUT_DIR, "*.pdf") # plus validation diagnostics (see abc library documentation)
#> [1] "validation_diagnostics.pdf"
validator$selected_simulations # top 3 models (stable abundance and high harvest)
#>  index  weight
#> 1     1 0.0346306
#> 2     4 0.0000000
#> 3     12 0.9779548

```

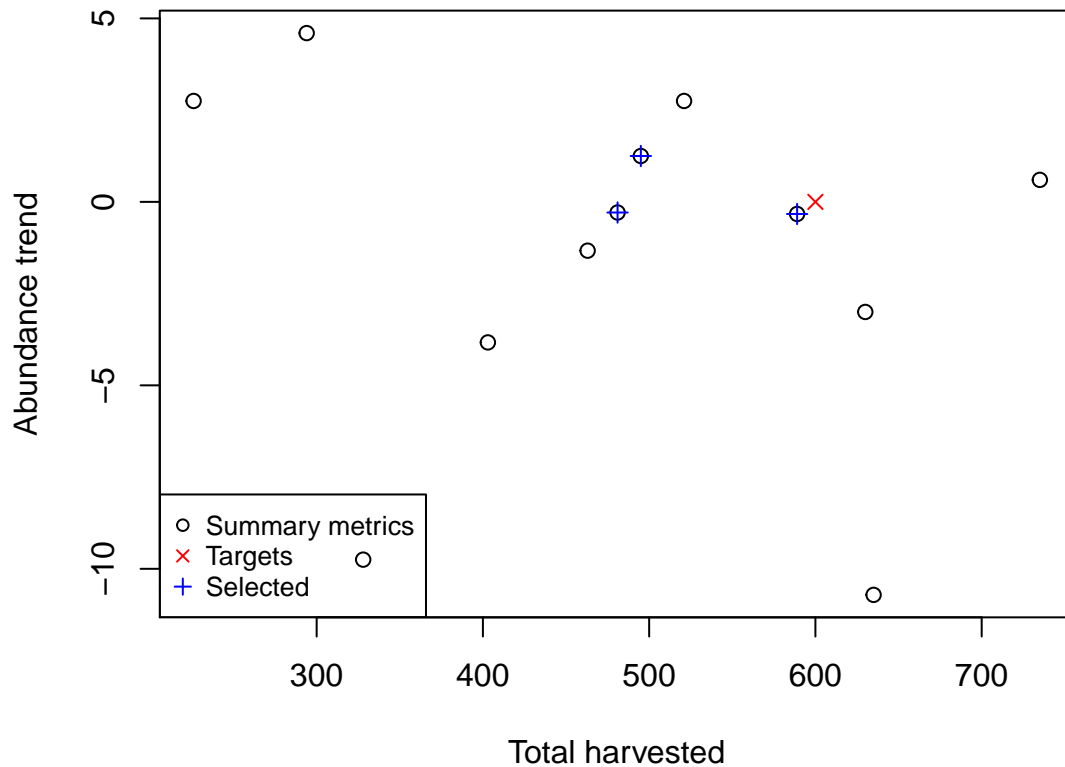
We encourage the user to examine the generated diagnostics (PDF) output file, and to become acquainted with the analysis that this information facilitates (see the *abc* documentation). As our simple example only uses two metrics/targets, we can visualize the congruence of the selected models with the targets via a simple plot.

```

# Plot the simulation, targets, and selected metrics
graphics::plot(x = results_manager$summary_metric_data$total_h,
              y = results_manager$summary_metric_data$trend_n,
              main = "Example model validation",
              xlab = "Total harvested", ylab = "Abundance trend")
graphics::points(x = 600, y = 0, col = "red", pch = 4)
selected_indices <- validator$selected_simulations$index
graphics::points(x = results_manager$summary_metric_data$total_h[selected_indices],
                y = results_manager$summary_metric_data$trend_n[selected_indices],
                col = "blue", pch = 3)
graphics::legend("bottomleft", legend = c("Summary metrics", "Targets", "Selected"),
                col = c(1, "red", "blue"), pch = c(1, 4, 3), cex = 0.8)

```


Example model validation



Summary

This demonstration has provided an overview of the *poems* workflow and modules via a simple population model example. We hope it has given you the foundation to progress to our more advanced Tasmanian *Thylacine* vignette, and towards utilizing the package for your own modeling projects.

Thank you :-)

References

- Beaumont, M. A., Zhang, W., & Balding, D. J. (2002). 'Approximate Bayesian computation in population genetics'. *Genetics*, vol. 162, no. 4, pp, 2025–2035.
- Csillery, K., Lemaire L., Francois O., & Blum M. (2015). 'abc: Tools for Approximate Bayesian Computation (ABC)'. *R package version 2.1*. Retrieved from <https://CRAN.R-project.org/package=abc>
- Grimm, V., Revilla, E., Berger, U., Jeltsch, F., Mooij, W. M., Railsback, S. F., Thulke, H. H., Weiner, J., Wiegand, T., DeAngelis, D. L., (2005). 'Pattern-Oriented Modeling of Agent-Based Complex Systems: Lessons from Ecology'. *Science* vol. 310, no. 5750, pp. 987–991.