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# Composite Indicator Analysis and Optimization (CIAO) Tool

*Codebook for practitioners, v.2*

**v.2: Revision January 2021**

(v.1: November 2018)

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## Program description

The Composite Indicator Analysis and Optimization (CIAO) tool is an expansion of the automated menu-version of the Matlab® toolbox presented by Becker (2017) for the advanced assessment of Composite Indicators (CIs). It was developed in connection to assessing the implicit weights of a novel index – the Electricity Supply Resilience Index (ESRI), developed within the Future Resilient Systems (FRS) program, at the Singapore-ETH Centre (SEC); see Gasser et al. (2017) and Lindén (2018) for further details. The scripts for the CIAO tool are written in MathWorks’s (<http://www.mathworks.com>) commercial software language, Matlab® 2017 version<sup>1</sup>, and contain five main steps: (1) a correlation analysis, (2) a weighting and aggregation of the indicators, (3) an estimation of the influence of each indicator (Si), (4) a decomposition of such influence, and (5) an optimization of the indicators’ weights.

Thus, the CIAO tool allows users to:

1. Perform a detailed examination of the linear and nonlinear relationships among (i) the set of indicators and (ii) between the indicators and the Composite Indicator (CI).
2. Assess how the indicators are “balanced” within the CI, i.e. to what extent the influence of each indicator matches its assigned weight.
3. Tune the weights of the indicators so that the influence of each one matches pre-defined importance values.

Details on the methodology are presented in Becker et al. (2017) and Lindén (2018).

Please NOTE that two additional commercial toolboxes are required to run the full analysis: The Statistics and Machine Learning Toolbox<sup>2</sup>, and the Optimization Toolbox<sup>3</sup>. Also, this package uses one freely available script from the MATLAB Central/File Exchange: *Random Vectors with Fixed Sum*<sup>4</sup>. For this analysis, it is used for generating random starting points for the weight-optimization.

This codebook presents the structure of the package to run the CIAO tool and provides details on each script.

*Revision January 2021 (v.2):*

Adaptations have been made to the CIAO code and the codebook to include the application of the modules to perform the analysis of information transfer. This includes the addition of a second objective for the weight optimization, i.e., maximize information transfer. The user can now choose between two objectives: (i) **Adjust** the contribution of each indicator in relation to its assigned weight, or (ii) **Maximize** the information transferred from the set of indicators to the CI.

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<sup>1</sup> Please NOTE that the current CIAO code was developed with Matlab® 2017 version. It is thus recommended not to run the CIAO code with another version of Matlab® as this might result in loss of functionality due to possible changes/updates of functions and underlying math libraries (implemented by MathWorks between Matlab® version updates), which the authors cannot take responsibility for.

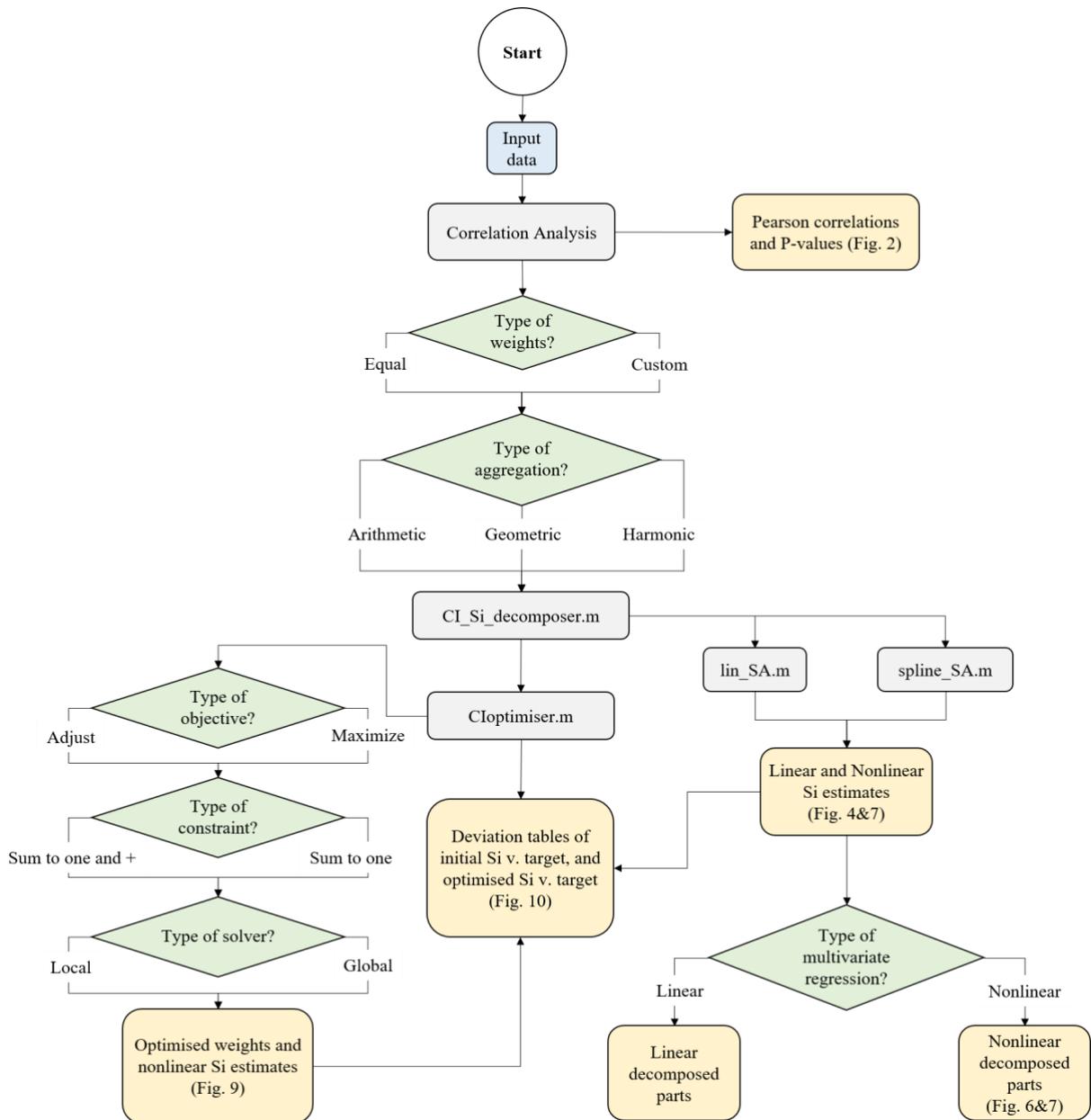
<sup>2</sup> URL: <https://se.mathworks.com/products/statistics.html>

<sup>3</sup> URL: <https://se.mathworks.com/products/optimization.html>

<sup>4</sup> Author: Roger Stafford, 2006 - URL: <https://se.mathworks.com/matlabcentral/fileexchange/9700-random-vectors-with-fixed-sum>

## Scripts

A flowchart of the CIAO tool is shown in Fig. 1. The scripts run in the tool are described in the following sections.



**Fig. 1.** Flow of the menu-version for implicit weights analysis and optimization in CIs; color coded accordingly: input (blue), outputs (orange), functions (grey), menu choices (green).

### *MENU\_version.m*

*MENU\_version.m* is the main script of this package. By running this script, a user-friendly GUI menu is initiated. Although most sub-function settings are automated, the GUI allows the user to “click” through the analysis with a certain amount of freedom to choose different settings (see the green menu choices in Fig. 1). For an explanation on what the script needs to run, see “Input” section at page 9. Default setting loads *test\_data.m*, which is a 58-by-4 matrix (see Input section). Then, the main script automatically generates a correlation matrix (with Pearson correlation coefficients) and its corresponding p-values for the matrix of indicators (Fig. 2).

```

17 %% Check correlation and P-values
18
19 [R,P] = corrcoef(X);
20 fprintf('Correlation matrix >>\n')
21 corr=array2table(round(R,4),'RowNames',rownamez, 'VariableNames', varnamez);
22 disp(corr)
23 fprintf('P-values >>\n')
24 pval=array2table(round(P,4),'RowNames',rownamez, 'VariableNames', varnamez);
25 disp(pval)

```

Command Window

```

Correlation matrix >>
      X_1      X_2      X_3      X_4
      ---      ---      ---      ---
X_1      1      0.6891  0.7286  0.4145
X_2  0.6891      1      0.8224  0.6249
X_3  0.7286  0.8224      1      0.5997
X_4  0.4145  0.6249  0.5997      1

P-values >>
      X_1      X_2      X_3      X_4
      ---      ---      ---      ---
X_1      1      0      0      0.0012
X_2      0      1      0      0
X_3      0      0      1      0
X_4  0.0012      0      0      1

```

Fig. 2. Output from the correlation analysis, which is obtained automatically by running the script. Note that default setting is to name variables  $X_i$ , where  $i=1,2,\dots,n$ .

Subsequently, the user is provided with the choice of what weights to apply to the indicators (either equal or custom) and also asked to choose an aggregation function (either arithmetic “ArAv”, geometric “GeAv”, or harmonic average “HarAv”); see Fig. 3. NOTE: The showcased results are obtained using the default data (*test\_data.m*), aggregated with arithmetic average and equal weights.

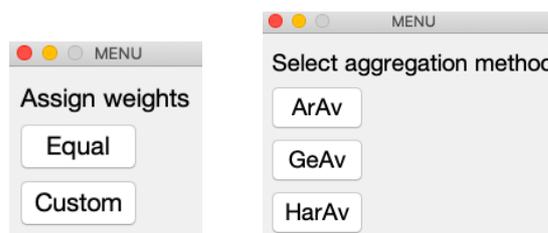


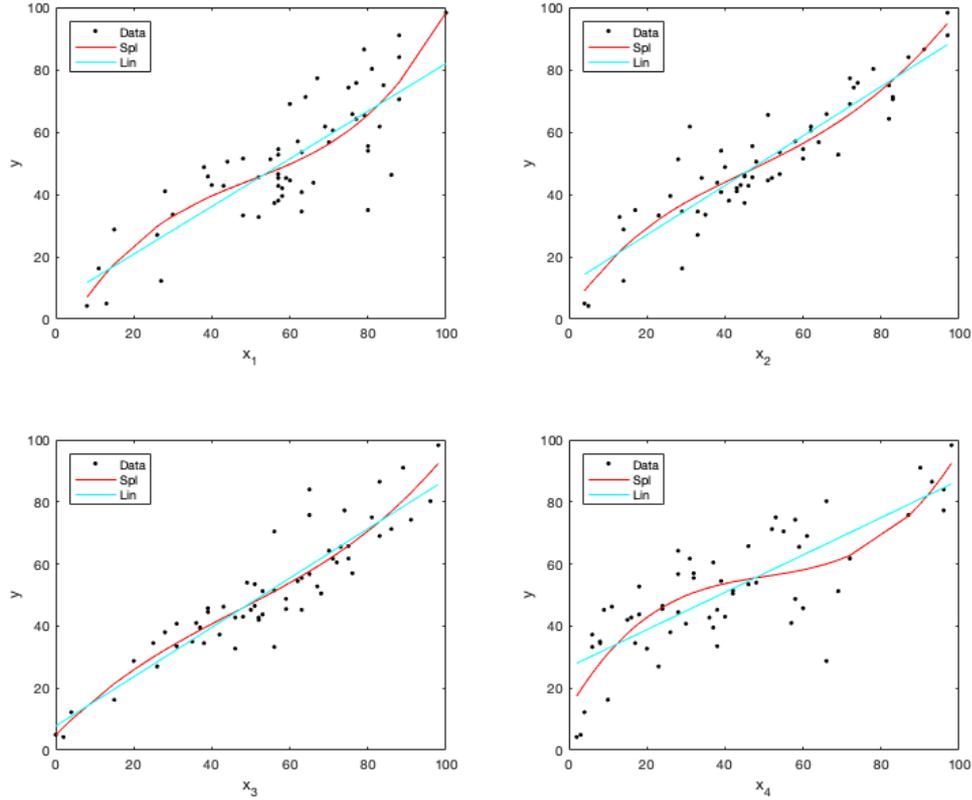
Fig. 3. The GUI menu for the selection of weights (left) and the selection of aggregation function (right).

### *CI\_Si\_decomposer.m*

*CI\_Si\_decomposer.m* is a function called by the main script *MENU\_version.m*. It uses the  $X$  and  $y$  data to provide the first order sensitivity indices (i.e. influence of each indicator,  $S_i$ ), plus their decomposed values into correlated and uncorrelated contributions. The function runs accordingly:

$$\text{function outSi} = \text{CI\_Si\_decomposer}(X, y, \text{plott})$$

The input parameters are the matrix of indicators ( $X$ ) and the vector ( $y$ ), which is the aggregated CI. Additionally, the choice of plotting the regression models ( $\text{plott}=1$ ) or not ( $\text{plott}=0$ ) can be made;  $\text{plott}=1$  gives scatter plots with splines and linear regression fits (Fig. 4).



**Fig. 4.** Output from the regression analysis. Plots showing linear (cyan) and the nonlinear splines (red) regression models.

By running this function, three steps are performed:

First, the function calls two sub-functions for the estimation of the first order sensitivity indices (Si):

*function [R2, par] = lin\_SA(y, X, plott)*

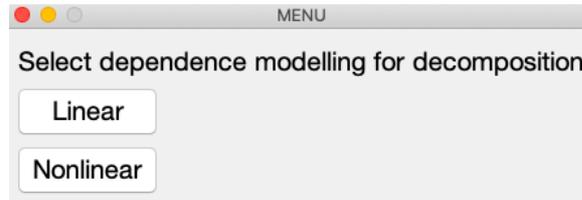
this calculates  $R^2$  (R2), i.e. the linear estimates of Si, for the X and y data. The linear regression parameters (e.g. the predicted outcome variables) are stored in a structured array (*par*) with two fields, including: the predicted outcome values of the regression model (*par.yhats*), and X-values sorted in ascending order (*par.xsrts*). Also,

*function [S\_E, S\_V, par] = spline\_SA(y, X, plott, p, knots)*

this uses the X and y data to calculate the Si, using nonlinear (splines) regression. The *CI\_Si\_decomposer.m* uses a cubic order of the spline ( $p=3$ ) and a row vector of knot numbers to try ( $knots = [1\ 2\ 3\ 4\ 5]$ ) as standard, but this can be altered if desired. The outputs are a column vector of first order sensitivity indices estimated via  $E(y|X_i)$  (S\_E) or  $V(y|X_i)$  (S\_V). The splines regression parameters are stored in a structured array (*par*) with eight fields, including: order of spline (*par.p*), trial knots (*par.trialknots*), the predicted outcome and variance values of the regression model (*par.yhats* & *par.varhats*), number of knots (*par.knotN*), derivatives (*par.derivs* & *par.derivs\_xx*), and X-values sorted in ascending order (*par.xsrts*).

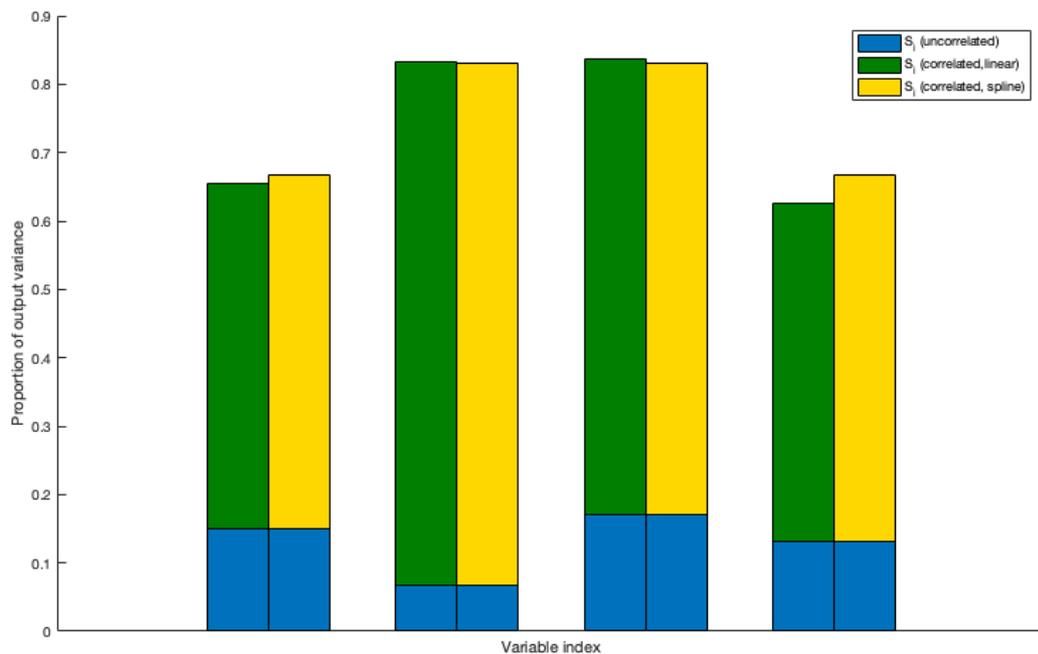
Second, a multivariate regression of  $X_i$  on  $X_{-i}$  (representing the matrix X without the  $i$ th column, i.e. indicator) is carried out, using either a standard linear model, according to the method of Xu and Gertner (2008), or MATLAB's inbuilt Gaussian process regression model

(*fitrgp*)<sup>5</sup> for a nonlinear model. The user is free to choose between these two models, as shown in Fig. 5.



**Fig. 5.** The GUI for the selection of model to perform the multivariate regression.

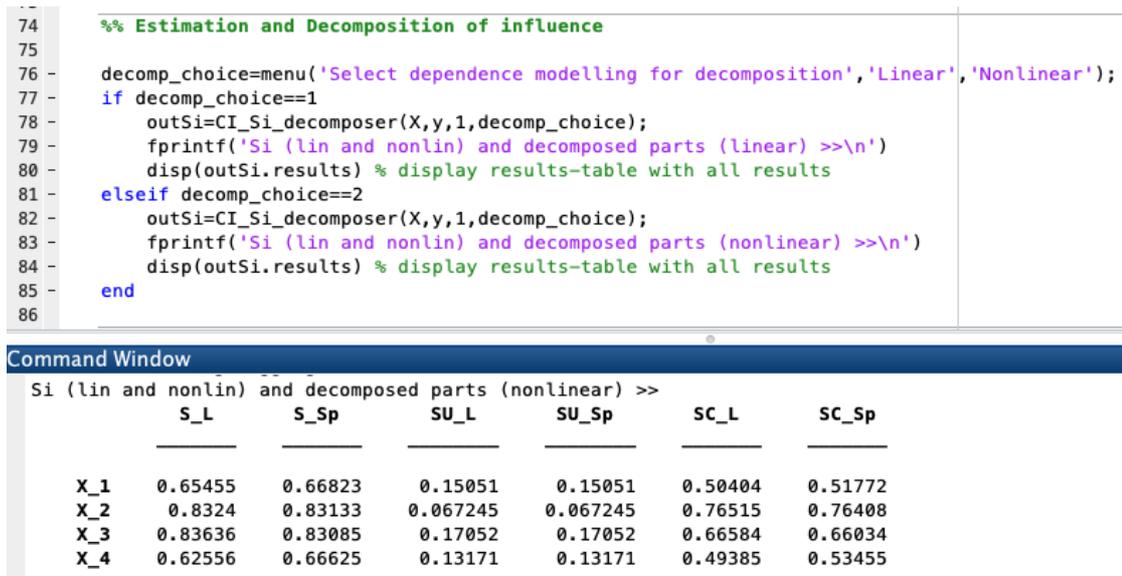
The results from the multivariate regression are then used to obtain a residual (i.e. what is “left” after correlation is removed). Subsequently, a nonlinear regression (splines) of  $y$  on the residuals of previous regression is performed to obtain the uncorrelated part (SU). Finally, the correlated part (SC) is obtained simply by subtracting the uncorrelated from the  $S_i$ , i.e.  $SC=S_i-SU$ . These are the decomposed values, which are then plotted in a bar chart (Fig. 6).



**Fig. 6.** The decomposed  $S_i$  values; full bars represent the  $S_i$  values estimated in the previous regression analysis. The uncorrelated parts (blue), estimated with nonlinear multivariate regression, and the resulting correlated parts, from either the linear (green) and nonlinear (yellow) regression results.

After these steps have been completed, the output of the *CI\_Si\_decomposer.m* is, beside the two plots, a structure array (*outSi*) with 3 fields containing: (1) linear regression results (*outSi.lin*), (2) spline regression results (*outSi.spline*), and (3) a summarized table of both estimates and the decomposed values (*outSi.results*), shown in Fig. 7.

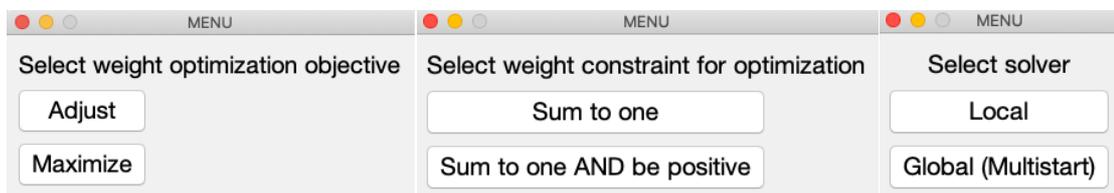
<sup>5</sup> NOTE: for this nonlinear dependence modelling, the Statistics and Machine Learning toolbox is required. If this is not detected, linear dependence modelling will be used.



**Fig. 7.** Output from the estimation and decomposition of influence. These are the values used for the bar chart in Fig. 6 above. S\_L and S\_Sp are the full bar values, i.e. the influence (Si), estimated with linear or nonlinear (splines) regression. The SU and SC are the uncorrelated and correlated parts, respectively.

### *CIoptimiser.m*

*CIoptimiser.m* is a self-contained function which optimizes CI weights, using nonlinear regression (penalized splines). It requires three choices (Fig. 8) prior to initiating the optimization algorithm<sup>6</sup>.



**Fig. 8.** The GUI for the choice of objective (left), optimization constraint (middle) and solver (right).

First, the objective of the optimization must be specified. There are two choices: (i) *Adjust* the contribution of each indicator in relation to its assigned weight, or (ii) *Maximize* the information transferred from the set of indicators to the CI. Second, a constraint on the weights must be chosen. The weights are either only constrained to sum to one (meaning they can take on negative values), OR they must sum to one but also be positive. Third, the choice is between two different types of solver, either a local or a global one. Although both are based on the same search method, the Nelder-Mead simplex, they differ in two aspects: (a) starting point(s) and (b) number of runs. The local solver uses the input weights as starting point, running the search algorithm once. Conversely, the global solver (MultiStart<sup>7</sup>) offers an option to specify the quantity of randomly generated starting points, re-running the search method for each one. This might be desired if the local solver fails to converge to an optimal solution<sup>8</sup>. However, for most cases, the local solver is sufficient and is thus recommended. After these choices are made, the function runs accordingly:

<sup>6</sup> NOTE: for constrained optimization, Optimization Toolbox is required. If this is not detected, will automatically revert to unconstrained.

<sup>7</sup> For a more detailed overview on this solver, see <https://se.mathworks.com/help/gads/multistart.html>

<sup>8</sup> For more info on this topic, see <https://se.mathworks.com/help/optim/ug/local-vs-global-optima.html>

*function outOpt = CIOptimiser(X, Sd, objective, constr, CLeq)*

this function uses the X data to find the optimal set of weights according to a predefined “target” (*Sd*), determined by the objective (*objective*) and a constraint defined by (*constr*); *constr=1* represents the sum-to-one constraint, and *constr=2* represents the sum-to-one constraint together with a positive value constraint. *CLeq* is the aggregation form (e.g. ‘ArAv’ for arithmetic) of the CI and it is set automatically by the *MENU\_version.m* script according to which function was chosen for aggregation at the beginning (Fig. 9). The output of this function is a structured array (*outOpt*) with six fields containing, for example: the optimized weights (*outOpt.wopt*), *Si* at optimized weights (*outOpt.Sopt*) and some other solver-specific parameters.

```

81  %% Optimization
82
83  w_choice=menu('Select weight constraint for optimization','Sum to one','Sum to one AND be positive');
84  if aggregation==1
85      outOpt=CIoptimiser(X,w,w_choice,'ArAv');
86  elseif aggregation==2
87      outOpt=CIoptimiser(X,w,w_choice,'GeAv');
88  elseif aggregation==3
89      outOpt=CIoptimiser(X,w,w_choice,'HarAv');

```

---

Command Window

Local minimum found that satisfies the constraints.

Optimization completed because the objective function is non-decreasing in feasible directions, to within the default value of the optimality tolerance, and constraints are satisfied to within the default value of the constraint tolerance.

<stopping criteria details>

Optimised weights >>

```

0.4443
0.0782
0.1002
0.3773

```

*Si* at optimised weights >>

```

0.7207
0.7207
0.7207
0.7207

```

**Fig. 9.** Output from weight-optimization, using the sum to one AND positive constraint. As seen from the stopping criterion, a local solver is used. Judging by the results, it successfully finds an optimal set of weights that achieves an equal *Si*.

- After all the steps are carried out, the main script *MENU\_version.m* automatically compiles two summarizing tables (Fig. 10), which showcase the discrepancy between initial *Si* values and the target values (weights), and optimized *Si* values and the target values (weights). For this purpose, both sets of *Si* values are first normalized to render them comparable to the weights (summing to one). Note that this is only performed if “Adjust” objective is selected.

```

Command Window

```

Normalised <i>Si</i> at initial weights >>				
	<i>Sis_norm</i>	<i>Sis_target</i>	<i>Sis_dev</i>	<i>Dev_ratio</i>
X_1	0.22299	0.25	-0.02701	'-11%'
X_2	0.27742	0.25	0.02742	' 11%'
X_3	0.27726	0.25	0.02726	' 11%'
X_4	0.22233	0.25	-0.02767	'-11%'

Normalised <i>Si</i> at optimised weights >>				
	<i>Sisopt_norm</i>	<i>Sis_target</i>	<i>Sisopt_dev</i>	<i>Dev_ratio</i>
X_1	0.25	0.25	0	'0%'
X_2	0.25	0.25	0	'0%'
X_3	0.25	0.25	0	'0%'
X_4	0.25	0.25	0	'0%'

**Fig. 10.** Summarizing tables of the results, comparing both the *Si* values at initial weights (above) and *Si* values at optimized weights (below) with reference to the target values (i.e. weights), in this case equal.

## Input

A matrix ( $X$ ), with  $N$ -alternatives (rows) by  $k$ -variables (columns). Default setting loads *test\_data.m*, which is a 58-by-4 matrix (58 alternatives, 4 indicators) containing data that is normalized to a 0-100 scale. To add your own dataset from, for example, an excel spreadsheet, simply use MATLAB's inbuilt import tool<sup>9</sup>. Follow the checklist below to confirm that data is imported correctly (format, name, etc.).

Checklist:

- 1) First, create/prepare an excel-file (.xlsx) with the input data. NOTE: The data must be normalized<sup>10</sup> prior to importing it. Example of a correct dataset can be found [here](#).
- 2) When importing the data, set the “output type” to be a numeric matrix (Fig. 11).

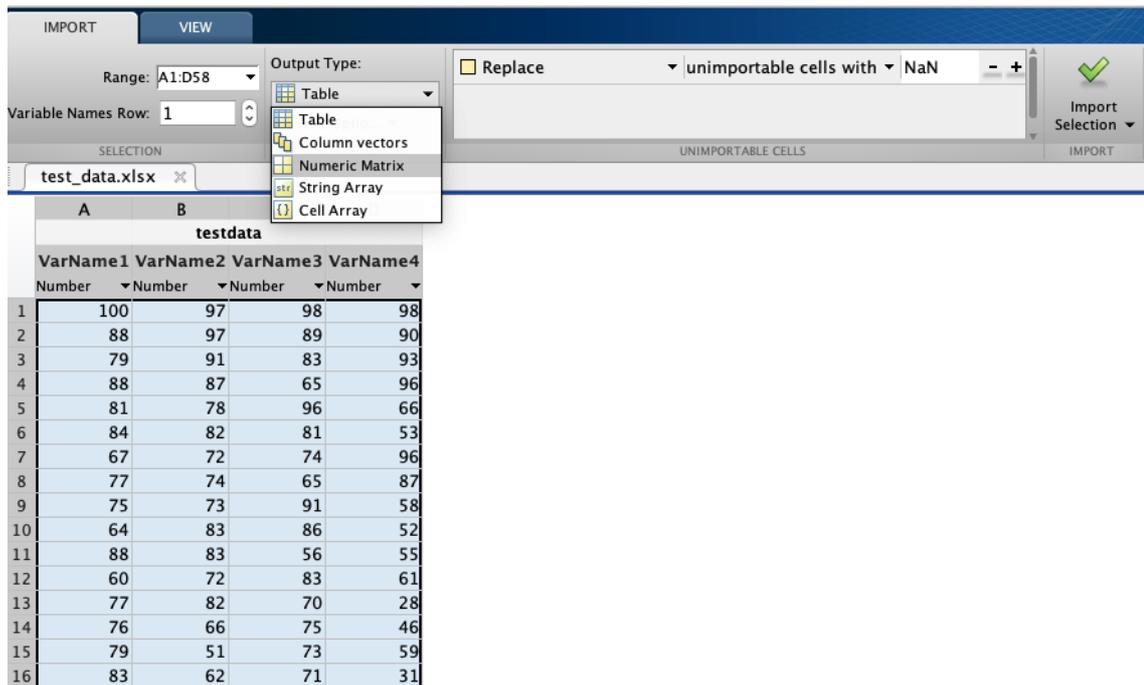
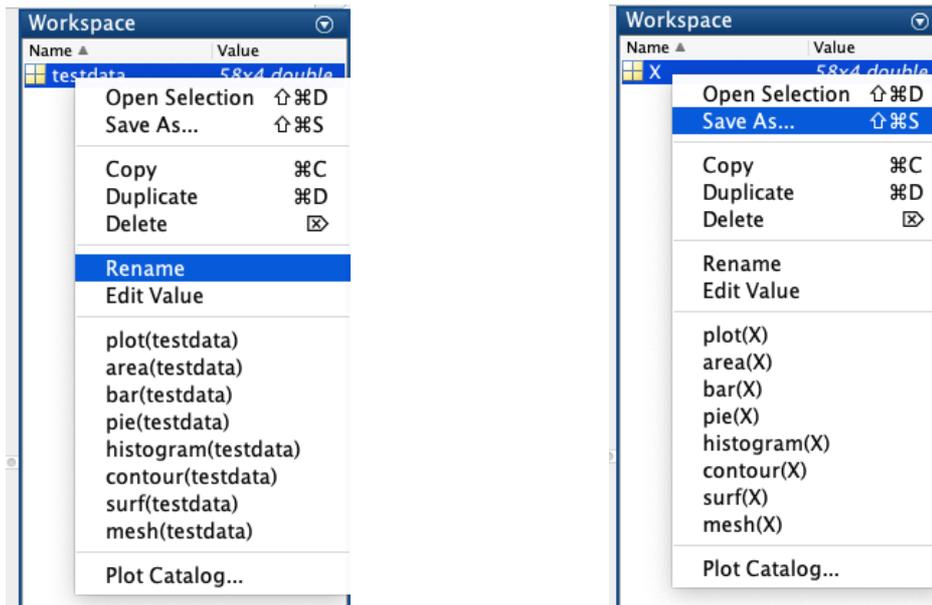


Fig. 11. Screenshot representing the choice of output types. Set to Numeric Matrix.

- 3) Be sure to assign your imported matrix to the global variable “X”, as this is fundamental for running the scripts. This is easily done by right-clicking on the matrix name in the workspace and selecting “Rename” (Fig. 12).
- 4) Save your imported X-matrix as a .mat file in the same folder as the scripts (Fig. 12).

<sup>9</sup> For more info on the import tool in MATLAB see for example, <https://se.mathworks.com/help/matlab/ref/importtool-app.html>

<sup>10</sup> Examples of appropriate normalization methods, (a) Min-Max (b) Target (c) Ranking. Avoid negative values. If standardization is used, it requires that all values are brought to a positive scale. This can be done by, for example, finding the largest negative value in the data set and adding its absolute form to all values (e.g.  $X + \text{abs}(\min(X))$ ). This ensures that the minimum value is zero. For more details on normalization methods, see OECD (2008).



**Fig. 12.** Screenshots representing the renaming of the imported matrix (left). Then, when renamed/assigned to the global variable X (right), select Save As... and save as a .mat file.

- 5) In the *MENU\_version.m* script, simply replace the “load test\_data” (Fig. 13) with the name that your data is saved as.

---

```

%% Load input data

load test_data % load X matrix containing indicator data

```

---

**Fig. 13.** The line in the script *MENU\_version.m* that loads the X matrix for analysis.

- 6) Now, everything should be set for running the analysis on your data.

## Output

Pearson correlations, linear and nonlinear regression estimates (Si), decomposed values (SU and SC), optimized Si and weights (wopt), and discrepancy tables.

## Acknowledgments

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