



## C interfaces to GALAHAD LSQP

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# Chapter 1

## GALAHAD C package Isqp

### 1.1 Introduction

#### 1.1.1 Purpose

This package uses a primal-dual interior-point trust-region method to solve the **linear** or **separable convex quadratic programming problem**

$$\text{minimize } \frac{1}{2} \sum_{j=1}^n w_j^2 (x_j - x_j^0)^2 + g^T x + f$$

subject to the general linear constraints

$$c_i^l \leq a_i^T x \leq c_i^u, \quad i = 1, \dots, m,$$

and the simple bound constraints

$$x_j^l \leq x_j \leq x_j^u, \quad j = 1, \dots, n,$$

where the vectors  $g$ ,  $w$ ,  $x^0$ ,  $c^l$ ,  $c^u$ ,  $x^l$ ,  $x^u$  and the scalar  $f$  are given. Any of the constraint bounds  $c_i^l$ ,  $c_i^u$ ,  $x_j^l$  and  $x_j^u$  may be infinite. Full advantage is taken of any zero coefficients in the matrix  $A$  of vectors  $a_i$ .

In the special case where  $w = 0$ ,  $g = 0$  and  $f = 0$ , the so-called analytic center of the feasible set will be found, while linear programming, or constrained least distance, problems may be solved by picking  $w = 0$ , or  $g = 0$  and  $f = 0$ , respectively.

The more-modern GALAHAD package CQP offers similar functionality, and is often to be preferred.

#### 1.1.2 Authors

N. I. M. Gould, STFC-Rutherford Appleton Laboratory, England, and Philippe L. Toint, University of Namur, Belgium.

C interface, additionally J. Fowkes, STFC-Rutherford Appleton Laboratory.

Julia interface, additionally A. Montoison and D. Orban, Polytechnique Montréal.

#### 1.1.3 Originally released

October 2001, C interface January 2022.

### 1.1.4 Terminology

The required solution  $x$  necessarily satisfies the primal optimality conditions

$$(1a) \quad Ax = c$$

and

$$(1b) \quad c^l \leq c \leq c^u, \quad x^l \leq x \leq x^u,$$

the dual optimality conditions

$$(2a) \quad W^2(x - x^0) + g = A^T y + z$$

where

$$(2b) \quad y = y^l + y^u, \quad z = z^l + z^u, \quad y^l \geq 0, \quad y^u \leq 0, \quad z^l \geq 0 \quad \text{and} \quad z^u \leq 0,$$

and the complementary slackness conditions

$$(3) \quad (Ax - c^l)^T y^l = 0, \quad (Ax - c^u)^T y^u = 0, \quad (x - x^l)^T z^l = 0 \quad \text{and} \quad (x - x^u)^T z^u = 0,$$

where the diagonal matrix  $W^2$  has diagonal entries  $w_j^2$ ,  $j = 1, \dots, n$ , where the vectors  $y$  and  $z$  are known as the Lagrange multipliers for the general linear constraints, and the dual variables for the bounds, respectively, and where the vector inequalities hold component-wise.

### 1.1.5 Method

Primal-dual interior point methods iterate towards a point that satisfies these conditions by ultimately aiming to satisfy (1a), (2a) and (3), while ensuring that (1b) and (2b) are satisfied as strict inequalities at each stage. Appropriate norms of the amounts by which (1a), (2a) and (3) fail to be satisfied are known as the primal and dual infeasibility, and the violation of complementary slackness, respectively. The fact that (1b) and (2b) are satisfied as strict inequalities gives such methods their other title, namely interior-point methods.

When  $w \neq 0$  or  $g \neq 0$ , the method aims at each stage to reduce the overall violation of (1a), (2a) and (3), rather than reducing each of the terms individually. Given an estimate  $v = (x, c, y, y^l, y^u, z, z^l, z^u)$  of the primal-dual variables, a correction  $\Delta v = \Delta(x, c, y, y^l, y^u, z, z^l, z^u)$  is obtained by solving a suitable linear system of Newton equations for the nonlinear systems (1a), (2a) and a parameterized "residual trajectory" perturbation of (3). An improved estimate  $v + \alpha \Delta v$  is then used, where the step-size  $\alpha$  is chosen as close to 1.0 as possible while ensuring both that (1b) and (2b) continue to hold and that the individual components which make up the complementary slackness (3) do not deviate too significantly from their average value. The parameter that controls the perturbation of (3) is ultimately driven to zero.

The Newton equations are solved by applying the GALAHAD matrix factorization package SBLS, but there are options to factorize the matrix as a whole (the so-called "augmented system" approach), to perform a block elimination first (the "Schur-complement" approach), or to let the method itself decide which of the two previous options is more appropriate. The "Schur-complement" approach is usually to be preferred when all the weights are nonzero or when every variable is bounded (at least one side), but may be inefficient if any of the columns of  $A$  is too dense.

When  $w = 0$  and  $g = 0$ , the method aims instead firstly to find an interior primal feasible point, that is to ensure that (1a) is satisfied. Once this has been achieved, attention is switched to minimizing the potential function

$$\phi(x, c) = \sum_{i=1}^m \log(c_i - c_i^l) + \sum_{i=1}^m \log(c_i^u - c_i) + \sum_{j=1}^n \log(x_j - x_j^l) + \sum_{j=1}^n \log(x_j^u - x_j),$$

while ensuring that (1a) remain satisfied and that  $x$  and  $c$  are strictly interior points for (1b). The global minimizer of this minimization problem is known as the analytic center of the feasible region, and may be viewed as a feasible point that is as far from the boundary of the constraints as possible. Note that terms in the above summations corresponding to infinite bounds are ignored, and that equality constraints are treated specially. Appropriate "primal" Newton corrections are used to generate a sequence of improving points converging to the analytic center, while the iteration is stabilized by performing insearches along these corrections with respect to  $\phi(x, c)$ .

In order to make the solution as efficient as possible, the variables and constraints are reordered internally by the GALAHAD package QPP prior to solution. In particular, fixed variables, and free (unbounded on both sides) constraints are temporarily removed. Optionally, the problem may be pre-processed temporarily to eliminate dependent constraints using the GALAHAD package FDC. This may improve the performance of the subsequent iteration.

### 1.1.6 Reference

The basic algorithm is a generalisation of those of

Y. Zhang (1994), On the convergence of a class of infeasible interior-point methods for the horizontal linear complementarity problem, SIAM J. Optimization 4(1) 208-227,

with a number of enhancements described by

A. R. Conn, N. I. M. Gould, D. Orban and Ph. L. Toint (1999). A primal-dual trust-region algorithm for minimizing a non-convex function subject to general inequality and linear equality constraints. Mathematical Programming 87 215-249.

### 1.1.7 Call order

To solve a given problem, functions from the lsqp package must be called in the following order:

- `lsqp_initialize` - provide default control parameters and set up initial data structures
- `lsqp_read_specfile` (optional) - override control values by reading replacement values from a file
- `lsqp_import` - set up problem data structures and fixed values
- `lsqp_reset_control` (optional) - possibly change control parameters if a sequence of problems are being solved
- `lsqp_solve_qp` - solve the quadratic program
- `lsqp_information` (optional) - recover information about the solution and solution process
- `lsqp_terminate` - deallocate data structures

See Section 4.1 for examples of use.

### 1.1.8 Unsymmetric matrix storage formats

The unsymmetric  $m$  by  $n$  constraint matrix  $A$  may be presented and stored in a variety of convenient input formats.

Both C-style (0 based) and fortran-style (1-based) indexing is allowed. Choose `control.f_indexing` as `false` for C style and `true` for fortran style; the discussion below presumes C style, but add 1 to indices for the corresponding fortran version.

Wrappers will automatically convert between 0-based (C) and 1-based (fortran) array indexing, so may be used transparently from C. This conversion involves both time and memory overheads that may be avoided by supplying data that is already stored using 1-based indexing.

#### 1.1.8.1 Dense storage format

The matrix  $A$  is stored as a compact dense matrix by rows, that is, the values of the entries of each row in turn are stored in order within an appropriate real one-dimensional array. In this case, component  $n * i + j$  of the storage array `A_val` will hold the value  $A_{ij}$  for  $0 \leq i \leq m - 1$ ,  $0 \leq j \leq n - 1$ .

### 1.1.8.2 Sparse co-ordinate storage format

Only the nonzero entries of the matrices are stored. For the  $l$ -th entry,  $0 \leq l \leq ne - 1$ , of  $A$ , its row index  $i$ , column index  $j$  and value  $A_{ij}$ ,  $0 \leq i \leq m - 1$ ,  $0 \leq j \leq n - 1$ , are stored as the  $l$ -th components of the integer arrays  $A\_row$  and  $A\_col$  and real array  $A\_val$ , respectively, while the number of nonzeros is recorded as  $A\_ne = ne$ .

### 1.1.8.3 Sparse row-wise storage format

Again only the nonzero entries are stored, but this time they are ordered so that those in row  $i$  appear directly before those in row  $i+1$ . For the  $i$ -th row of  $A$  the  $i$ -th component of the integer array  $A\_ptr$  holds the position of the first entry in this row, while  $A\_ptr(m)$  holds the total number of entries. The column indices  $j$ ,  $0 \leq j \leq n - 1$ , and values  $A_{ij}$  of the nonzero entries in the  $i$ -th row are stored in components  $l = A\_ptr(i), \dots, A\_ptr(i+1)-1$ ,  $0 \leq i \leq m - 1$ , of the integer array  $A\_col$ , and real array  $A\_val$ , respectively. For sparse matrices, this scheme almost always requires less storage than its predecessor.



## Chapter 2

# File Index

### 2.1 File List

Here is a list of all files with brief descriptions:

<a href="#">galahad_lsqp.h</a> . . . . .	7
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## Chapter 3

# File Documentation

### 3.1 galahad\_lsqp.h File Reference

```
#include <stdbool.h>
#include <stdint.h>
#include "galahad_precision.h"
#include "galahad_cfunctions.h"
#include "galahad_fdc.h"
#include "galahad_sbfs.h"
```

#### Data Structures

- struct [lsqp\\_control\\_type](#)
- struct [lsqp\\_time\\_type](#)
- struct [lsqp\\_inform\\_type](#)

#### Functions

- void [lsqp\\_initialize](#) (void \*\*data, struct [lsqp\\_control\\_type](#) \*control, int \*status)
- void [lsqp\\_read\\_specfile](#) (struct [lsqp\\_control\\_type](#) \*control, const char specfile[])
- void [lsqp\\_import](#) (struct [lsqp\\_control\\_type](#) \*control, void \*\*data, int \*status, int n, int m, const char A\_type[], int A\_ne, const int A\_row[], const int A\_col[], const int A\_ptr[])
- void [lsqp\\_reset\\_control](#) (struct [lsqp\\_control\\_type](#) \*control, void \*\*data, int \*status)
- void [lsqp\\_solve\\_qp](#) (void \*\*data, int \*status, int n, int m, const real\_wp\_ w[], const real\_wp\_ x0[], const real\_wp\_ g[], const real\_wp\_ f, int a\_ne, const real\_wp\_ A\_val[], const real\_wp\_ c\_l[], const real\_wp\_ c\_u[], const real\_wp\_ x\_l[], const real\_wp\_ x\_u[], real\_wp\_ x[], real\_wp\_ c[], real\_wp\_ y[], real\_wp\_ z[], int x\_stat[], int c\_stat[])
- void [lsqp\\_information](#) (void \*\*data, struct [lsqp\\_inform\\_type](#) \*inform, int \*status)
- void [lsqp\\_terminate](#) (void \*\*data, struct [lsqp\\_control\\_type](#) \*control, struct [lsqp\\_inform\\_type](#) \*inform)

#### 3.1.1 Data Structure Documentation

##### 3.1.1.1 struct [lsqp\\_control\\_type](#)

control derived type as a C struct

##### Examples

[lsqpt.c](#), and [lsqptf.c](#).

## Data Fields

bool	f_indexing	use C or Fortran sparse matrix indexing
int	error	error and warning diagnostics occur on stream error
int	out	general output occurs on stream out
int	print_level	the level of output required is specified by print_level
int	start_print	any printing will start on this iteration
int	stop_print	any printing will stop on this iteration
int	maxit	at most maxit inner iterations are allowed
int	factor	the factorization to be used. Possible values are <ul style="list-style-type: none"> <li>• 0 automatic</li> <li>• 1 Schur-complement factorization</li> <li>• 2 augmented-system factorization</li> </ul>
int	max_col	the maximum number of nonzeros in a column of A which is permitted with the Schur-complement factorization
int	indmin	an initial guess as to the integer workspace required by SBLS
int	valmin	an initial guess as to the real workspace required by SBLS
int	itref_max	the maximum number of iterative refinements allowed
int	infeas_max	the number of iterations for which the overall infeasibility of the problem is not reduced by at least a factor .reduce_infeas before the problem is flagged as infeasible (see reduce_infeas)
int	muzero_fixed	the initial value of the barrier parameter will not be changed for the first muzero_fixed iterations
int	restore_problem	indicate whether and how much of the input problem should be restored on output. Possible values are <ul style="list-style-type: none"> <li>• 0 nothing restored</li> <li>• 1 scalar and vector parameters</li> <li>• 2 all parameters</li> </ul>

## Data Fields

int	indicator_type	<p>specifies the type of indicator function used. Possible values are</p> <ul style="list-style-type: none"> <li>• 1 primal indicator: constraint active if and only if the distance to nearest bound <math>\leq</math> .indicator_p_tol</li> <li>• 2 primal-dual indicator: constraint active if and only if the distance to nearest bound <math>\leq</math> .indicator_tol_pd * size of corresponding multiplier</li> <li>• 3 primal-dual indicator: constraint active if and only if the distance to the nearest bound <math>\leq</math> .indicator_tol_tapia * distance to same bound at previous iteration</li> </ul>
int	extrapolate	<p>should extrapolation be used to track the central path? Possible values</p> <ul style="list-style-type: none"> <li>• 0 never</li> <li>• 1 after the final major iteration</li> <li>• 2 at each major iteration (unused at present)</li> </ul>
int	path_history	the maximum number of previous path points to use when fitting the data (unused at present)
int	path_derivatives	the maximum order of path derivative to use (unused at present)
int	fit_order	the order of (Puisseux) series to fit to the path data: \$ to fit all data (unused at present)
int	sif_file_device	specifies the unit number to write generated SIF file describing the current problem
real_wp_	infinity	any bound larger than infinity in modulus will be regarded as infinite
real_wp_	stop_p	the required accuracy for the primal infeasibility
real_wp_	stop_d	the required accuracy for the dual infeasibility
real_wp_	stop_c	the required accuracy for the complementarity
real_wp_	prfeas	initial primal variables will not be closer than prfeas from their bounds
real_wp_	dufeas	initial dual variables will not be closer than dufears from their bounds
real_wp_	muzero	the initial value of the barrier parameter. If muzero is not positive, it will be reset to an appropriate value
real_wp_	reduce_infeas	if the overall infeasibility of the problem is not reduced by at least a factor reduce_infeas over .infeas_max iterations, the problem is flagged as infeasible (see infeas_max)

## Data Fields

real_wp_	potential_unbounded	if $W=0$ and the potential function value is smaller than <code>potential_unbounded * number of one-sided bounds</code> , the analytic center will be flagged as unbounded
real_wp_	pivot_tol	the threshold pivot used by the matrix factorization. See the documentation for SBLS for details
real_wp_	pivot_tol_for_dependencies	the threshold pivot used by the matrix factorization when attempting to detect linearly dependent constraints. See the documentation for SBLS for details
real_wp_	zero_pivot	any pivots smaller than <code>zero_pivot</code> in absolute value will be regarded to zero when attempting to detect linearly dependent constraints
real_wp_	identical_bounds_tol	any pair of constraint bounds ( $c_l, c_u$ ) or ( $x_l, x_u$ ) that are closer than <code>identical_bounds_tol</code> will be reset to the average of their values
real_wp_	mu_min	start terminal extrapolation when $\mu$ reaches <code>mu_min</code>
real_wp_	indicator_tol_p	if <code>.indicator_type = 1</code> , a constraint/bound will be deemed to be active if and only if the distance to nearest bound $\leq$ <code>.indicator_p_tol</code>
real_wp_	indicator_tol_pd	if <code>.indicator_type = 2</code> , a constraint/bound will be deemed to be active if and only if the distance to nearest bound $\leq$ <code>.indicator_tol_pd * size of corresponding multiplier</code>
real_wp_	indicator_tol_tapia	if <code>.indicator_type = 3</code> , a constraint/bound will be deemed to be active if and only if the distance to nearest bound $\leq$ <code>.indicator_tol_tapia * distance to same bound at previous iteration</code>
real_wp_	cpu_time_limit	the maximum CPU time allowed (-ve means infinite)
real_wp_	clock_time_limit	the maximum elapsed clock time allowed (-ve means infinite)
bool	remove_dependencies	the equality constraints will be preprocessed to remove any linear dependencies if true
bool	treat_zero_bounds_as_general	any problem bound with the value zero will be treated as if it were a general value if true
bool	just_feasible	if <code>.just_feasible</code> is true, the algorithm will stop as soon as a feasible point is found. Otherwise, the optimal solution to the problem will be found
bool	getdua	if <code>.getdua</code> is true, advanced initial values are obtained for the dual variables
bool	puiseux	If extrapolation is to be used, decide between Puiseux and Taylor series.
bool	feasol	if <code>.feasol</code> is true, the final solution obtained will be perturbed so that variables close to their bounds are moved onto these bounds

## Data Fields

bool	balance_initial_complentarity	if .balance_initial_complentarity is true, the initial complemetarity is required to be balanced
bool	use_corrector	if .use_corrector, a corrector step will be used
bool	array_syntax_worse_than_do_loop	if .array_syntax_worse_than_do_loop is true, f77-style do loops will be used rather than f90-style array syntax for vector operations
bool	space_critical	if .space_critical true, every effort will be made to use as little space as possible. This may result in longer computation time
bool	deallocate_error_fatal	if .deallocate_error_fatal is true, any array/pointer deallocation error will terminate execution. Otherwise, computation will continue
bool	generate_sif_file	if .generate_sif_file is .true. if a SIF file describing the current problem is to be generated
char	sif_file_name[31]	name of generated SIF file containing input problem
char	prefix[31]	all output lines will be prefixed by .prefix(2:LEN(TRIM(.prefix))-1) where .prefix contains the required string enclosed in quotes, e.g. "string" or 'string'
struct fdc_control_type	fdc_control	control parameters for FDC
struct sbls_control_type	sbls_control	control parameters for SBLS

## 3.1.1.2 struct lsqp\_time\_type

time derived type as a C struct

## Data Fields

real_wp_	total	the total CPU time spent in the package
real_wp_	preprocess	the CPU time spent preprocessing the problem
real_wp_	find_dependent	the CPU time spent detecting linear dependencies
real_wp_	analyse	the CPU time spent analysing the required matrices prior to factorization
real_wp_	factorize	the CPU time spent factorizing the required matrices
real_wp_	solve	the CPU time spent computing the search direction
real_wp_	clock_total	the total clock time spent in the package
real_wp_	clock_preprocess	the clock time spent preprocessing the problem
real_wp_	clock_find_dependent	the clock time spent detecting linear dependencies
real_wp_	clock_analyse	the clock time spent analysing the required matrices prior to factorization
real_wp_	clock_factorize	the clock time spent factorizing the required matrices
real_wp_	clock_solve	the clock time spent computing the search direction

### 3.1.1.3 struct lsqp\_inform\_type

inform derived type as a C struct

#### Examples

[lsqpt.c](#), and [lsqptf.c](#).

#### Data Fields

int	status	return status. See LSQP_solve for details
int	alloc_status	the status of the last attempted allocation/deallocation
char	bad_alloc[81]	the name of the array for which an allocation/deallocation error occurred
int	iter	the total number of iterations required
int	factorization_status	the return status from the factorization
int64_t	factorization_integer	the total integer workspace required for the factorization
int64_t	factorization_real	the total real workspace required for the factorization
int	nfacts	the total number of factorizations performed
int	nbacts	the total number of "wasted" function evaluations during the linesearch
real_wp_	obj	the value of the objective function at the best estimate of the solution determined by LSQP_solve_qp
real_wp_	potential	the value of the logarithmic potential function sum -log(distance to constraint boundary)
real_wp_	non_negligible_pivot	the smallest pivot which was not judged to be zero when detecting linear dependent constraints
bool	feasible	is the returned "solution" feasible?
struct <a href="#">lsqp_time_type</a>	time	timings (see above)
struct <a href="#">fdc_inform_type</a>	fdc_inform	inform parameters for FDC
struct <a href="#">sbls_inform_type</a>	sbls_inform	inform parameters for SBLS

## 3.1.2 Function Documentation

### 3.1.2.1 lsqp\_initialize()

```
void lsqp_initialize (
    void ** data,
    struct lsqp\_control\_type * control,
    int * status )
```

Set default control values and initialize private data

#### Parameters

in, out	<i>data</i>	holds private internal data
out	<i>control</i>	is a struct containing control information (see <a href="#">lsqp_control_type</a> )



## Parameters

out	<i>status</i>	is a scalar variable of type int, that gives the exit status from the package. Possible values are (currently): <ul style="list-style-type: none"> <li>• 0. The import was succesful.</li> </ul>
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## Examples

[lsqpt.c](#), and [lsqptf.c](#).

## 3.1.2.2 lsqp\_read\_specfile()

```
void lsqp_read_specfile (
    struct lsqp\_control\_type * control,
    const char specfile[] )
```

Read the content of a specification file, and assign values associated with given keywords to the corresponding control parameters. By default, the spcification file will be named RUNLSQP.SPC and lie in the current directory. Refer to Table 2.1 in the fortran documentation provided in \$GALAHAD/doc/lsqp.pdf for a list of keywords that may be set.

## Parameters

in, out	<i>control</i>	is a struct containing control information (see <a href="#">lsqp_control_type</a> )
in	<i>specfile</i>	is a character string containing the name of the specification file

## 3.1.2.3 lsqp\_import()

```
void lsqp_import (
    struct lsqp\_control\_type * control,
    void ** data,
    int * status,
    int n,
    int m,
    const char A_type[],
    int A_ne,
    const int A_row[],
    const int A_col[],
    const int A_ptr[] )
```

Import problem data into internal storage prior to solution.

## Parameters

in	<i>control</i>	is a struct whose members provide control paramters for the remaining pcedures (see <a href="#">lsqp_control_type</a> )
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## Parameters

in, out	<i>data</i>	holds private internal data
in, out	<i>status</i>	<p>is a scalar variable of type int, that gives the exit status from the package. Possible values are:</p> <ul style="list-style-type: none"> <li>• 0. The import was succesful</li> <li>• -1. An allocation error occurred. A message indicating the offending array is written on unit.control.error, and the returned allocation status and a string containing the name of the offending array are held in inform.alloc_status and inform.bad_alloc respectively.</li> <li>• -2. A deallocation error occurred. A message indicating the offending array is written on unit.control.error and the returned allocation status and a string containing the name of the offending array are held in inform.alloc_status and inform.bad_alloc respectively.</li> <li>• -3. The restrictions <math>n &gt; 0</math> or <math>m &gt; 0</math> or requirement that a type contains its relevant string 'dense', 'coordinate', 'sparse_by_rows', 'diagonal', 'scaled_identity', 'identity', 'zero' or 'none' has been violated.</li> </ul>
in	<i>n</i>	is a scalar variable of type int, that holds the number of variables.
in	<i>m</i>	is a scalar variable of type int, that holds the number of general linear constraints.
in	<i>A_type</i>	is a one-dimensional array of type char that specifies the <a href="#">unsymmetric storage scheme</a> used for the constraint Jacobian, <i>A</i> . It should be one of 'coordinate', 'sparse_by_rows' or 'dense'; lower or upper case variants are allowed.
in	<i>A_ne</i>	is a scalar variable of type int, that holds the number of entries in <i>A</i> in the sparse co-ordinate storage scheme. It need not be set for any of the other schemes.
in	<i>A_row</i>	is a one-dimensional array of size <i>A_ne</i> and type int, that holds the row indices of <i>A</i> in the sparse co-ordinate storage scheme. It need not be set for any of the other schemes, and in this case can be NULL.
in	<i>A_col</i>	is a one-dimensional array of size <i>A_ne</i> and type int, that holds the column indices of <i>A</i> in either the sparse co-ordinate, or the sparse row-wise storage scheme. It need not be set when the dense or diagonal storage schemes are used, and in this case can be NULL.
in	<i>A_ptr</i>	is a one-dimensional array of size $n+1$ and type int, that holds the starting position of each row of <i>A</i> , as well as the total number of entries, in the sparse row-wise storage scheme. It need not be set when the other schemes are used, and in this case can be NULL.

## Examples

[lsqpt.c](#), and [lsqptf.c](#).

## 3.1.2.4 lsqp\_reset\_control()

```
void lsqp_reset_control (
    struct lsqp_control_type * control,
    void ** data,
    int * status )
```

Reset control parameters after import if required.

## Parameters

in	<i>control</i>	is a struct whose members provide control paramters for the remaining prcedures (see <a href="#">lsqp_control_type</a> )
in, out	<i>data</i>	holds private internal data
in, out	<i>status</i>	is a scalar variable of type int, that gives the exit status from the package. Possible values are: <ul style="list-style-type: none"> <li>• 0. The import was succesful.</li> </ul>

## 3.1.2.5 lsqp\_solve\_qp()

```

void lsqp_solve_qp (
    void ** data,
    int * status,
    int n,
    int m,
    const real_wp_ w[],
    const real_wp_ x0[],
    const real_wp_ g[],
    const real_wp_ f,
    int a_ne,
    const real_wp_ A_val[],
    const real_wp_ c_l[],
    const real_wp_ c_u[],
    const real_wp_ x_l[],
    const real_wp_ x_u[],
    real_wp_ x[],
    real_wp_ c[],
    real_wp_ y[],
    real_wp_ z[],
    int x_stat[],
    int c_stat[] )

```

Solve the separable convex quadratic program.

## Parameters

in, out	<i>data</i>	holds private internal data
---------	-------------	-----------------------------

## Parameters

<code>in, out</code>	<code>status</code>	<p>is a scalar variable of type int, that gives the entry and exit status from the package. Possible exit are:</p> <ul style="list-style-type: none"> <li>• 0. The run was succesful</li> <li>• -1. An allocation error occurred. A message indicating the offending array is written on unit control.error, and the returned allocation status and a string containing the name of the offending array are held in <code>inform.alloc_status</code> and <code>inform.bad_alloc</code> respectively.</li> <li>• -2. A deallocation error occurred. A message indicating the offending array is written on unit control.error and the returned allocation status and a string containing the name of the offending array are held in <code>inform.alloc_status</code> and <code>inform.bad_alloc</code> respectively.</li> <li>• -3. The restrictions <math>n &gt; 0</math> and <math>m &gt; 0</math> or requirement that a type contains its relevant string 'dense', 'coordinate', 'sparse_by_rows', 'diagonal', 'scaled_identity', 'identity', 'zero' or 'none' has been violated.</li> <li>• -5. The simple-bound constraints are inconsistent.</li> <li>• -7. The constraints appear to have no feasible point.</li> <li>• -9. The analysis phase of the factorization failed; the return status from the factorization package is given in the component <code>inform.factor_status</code></li> <li>• -10. The factorization failed; the return status from the factorization package is given in the component <code>inform.factor_status</code>.</li> <li>• -11. The solution of a set of linear equations using factors from the factorization package failed; the return status from the factorization package is given in the component <code>inform.factor_status</code>.</li> <li>• -16. The problem is so ill-conditioned that further progress is impossible.</li> <li>• -17. The step is too small to make further impact.</li> <li>• -18. Too many iterations have been performed. This may happen if <code>control.maxit</code> is too small, but may also be symptomatic of a badly scaled problem.</li> <li>• -19. The CPU time limit has been reached. This may happen if <code>control.cpu_time_limit</code> is too small, but may also be symptomatic of a badly scaled problem.</li> </ul>
<code>in</code>	<code>n</code>	is a scalar variable of type int, that holds the number of variables
<code>in</code>	<code>m</code>	is a scalar variable of type int, that holds the number of general linear constraints.
<code>in</code>	<code>w</code>	is a one-dimensional array of size n and type double, that holds the values of the weights $w$ .
<code>in</code>	<code>x0</code>	is a one-dimensional array of size n and type double, that holds the values of the shifts $x^0$ .
<code>in</code>	<code>g</code>	is a one-dimensional array of size n and type double, that holds the linear term $g$ of the objective function. The j-th component of $g$ , $j = 0, \dots, n-1$ , contains $g_j$ .
<code>in</code>	<code>f</code>	is a scalar of type double, that holds the constant term $f$ of the objective function.
<code>in</code>	<code>a_ne</code>	is a scalar variable of type int, that holds the number of entries in the constraint Jacobian matrix $A$ .
<code>in</code>	<code>A_val</code>	is a one-dimensional array of size <code>a_ne</code> and type double, that holds the values of the entries of the constraint Jacobian matrix $A$ in any of the available storage schemes.
<code>in</code>	<code>c_l</code>	is a one-dimensional array of size m and type double, that holds the lower bounds $c^l$ on the constraints $Ax$ . The i-th component of <code>c_l</code> , $i = 0, \dots, m-1$ , contains $c_i^l$ .

## Parameters

in	<code>c_u</code>	is a one-dimensional array of size m and type double, that holds the upper bounds $c^l$ on the constraints $Ax$ . The i-th component of <code>c_u</code> , $i = 0, \dots, m-1$ , contains $c_i^u$ .
in	<code>x_l</code>	is a one-dimensional array of size n and type double, that holds the lower bounds $x^l$ on the variables $x$ . The j-th component of <code>x_l</code> , $j = 0, \dots, n-1$ , contains $x_j^l$ .
in	<code>x_u</code>	is a one-dimensional array of size n and type double, that holds the upper bounds $x^l$ on the variables $x$ . The j-th component of <code>x_u</code> , $j = 0, \dots, n-1$ , contains $x_j^l$ .
in, out	<code>x</code>	is a one-dimensional array of size n and type double, that holds the values $x$ of the optimization variables. The j-th component of <code>x</code> , $j = 0, \dots, n-1$ , contains $x_j$ .
out	<code>c</code>	is a one-dimensional array of size m and type double, that holds the residual $c(x)$ . The i-th component of <code>c</code> , $i = 0, \dots, m-1$ , contains $c_i(x)$ .
in, out	<code>y</code>	is a one-dimensional array of size n and type double, that holds the values $y$ of the Lagrange multipliers for the general linear constraints. The j-th component of <code>y</code> , $i = 0, \dots, m-1$ , contains $y_i$ .
in, out	<code>z</code>	is a one-dimensional array of size n and type double, that holds the values $z$ of the dual variables. The j-th component of <code>z</code> , $j = 0, \dots, n-1$ , contains $z_j$ .
out	<code>x_stat</code>	is a one-dimensional array of size n and type int, that gives the optimal status of the problem variables. If <code>x_stat(j)</code> is negative, the variable $x_j$ most likely lies on its lower bound, if it is positive, it lies on its upper bound, and if it is zero, it lies between its bounds.
out	<code>c_stat</code>	is a one-dimensional array of size m and type int, that gives the optimal status of the general linear constraints. If <code>c_stat(i)</code> is negative, the constraint value $a_i^T x$ most likely lies on its lower bound, if it is positive, it lies on its upper bound, and if it is zero, it lies between its bounds.

## Examples

[lsqpt.c](#), and [lsqptf.c](#).

## 3.1.2.6 lsqp\_information()

```
void lsqp_information (
    void ** data,
    struct lsqp_inform_type * inform,
    int * status )
```

Provides output information.

## Parameters

in, out	<code>data</code>	holds private internal data
out	<code>inform</code>	is a struct containing output information (see <a href="#">lsqp_inform_type</a> )
out	<code>status</code>	is a scalar variable of type int, that gives the exit status from the package. Possible values are (currently): <ul style="list-style-type: none"> <li>• 0. The values were recorded succesfully</li> </ul>

## Examples

[lsqpt.c](#), and [lsqptf.c](#).

## 3.1.2.7 lsqp\_terminate()

```
void lsqp_terminate (
    void ** data,
    struct lsqp_control_type * control,
    struct lsqp_inform_type * inform )
```

Deallocate all internal private storage.

## Parameters

in, out	<i>data</i>	holds private internal data
out	<i>control</i>	is a struct containing control information (see <a href="#">lsqp_control_type</a> )
out	<i>inform</i>	is a struct containing output information (see <a href="#">lsqp_inform_type</a> )

## Examples

[lsqpt.c](#), and [lsqptf.c](#).

## Chapter 4

# Example Documentation

### 4.1 lsqpt.c

This is an example of how to use the package to solve a quadratic program. A variety of supported Hessian and constraint matrix storage formats are shown.

Notice that C-style indexing is used, and that this is flagged by setting `control.f_indexing` to `false`.

```
/* lsqpt.c */
/* Full test for the LSQP C interface using C sparse matrix indexing */
#include <stdio.h>
#include <math.h>
#include "galahad_precision.h"
#include "galahad_cfunctions.h"
#include "galahad_lsqp.h"
int main(void) {
    // Derived types
    void *data;
    struct lsqp_control_type control;
    struct lsqp_inform_type inform;
    // Set problem data
    int n = 3; // dimension
    int m = 2; // number of general constraints
    real_wp_ g[] = {0.0, 2.0, 0.0}; // linear term in the objective
    real_wp_ f = 1.0; // constant term in the objective
    int A_ne = 4; // Jacobian elements
    int A_row[] = {0, 0, 1, 1}; // row indices
    int A_col[] = {0, 1, 1, 2}; // column indices
    int A_ptr[] = {0, 2, 4}; // row pointers
    real_wp_ A_val[] = {2.0, 1.0, 1.0, 1.0}; // values
    real_wp_ c_l[] = {1.0, 2.0}; // constraint lower bound
    real_wp_ c_u[] = {2.0, 2.0}; // constraint upper bound
    real_wp_ x_l[] = {-1.0, - INFINITY, - INFINITY}; // variable lower bound
    real_wp_ x_u[] = {1.0, INFINITY, 2.0}; // variable upper bound
    real_wp_ w[] = {1.0, 1.0, 1.0};
    real_wp_ x_0[] = {0.0, 0.0, 0.0};
    // Set output storage
    real_wp_ c[m]; // constraint values
    int x_stat[n]; // variable status
    int c_stat[m]; // constraint status
    char st;
    int status;
    printf(" C sparse matrix indexing\n\n");
    printf(" basic tests of qp storage formats\n\n");
    for( int d=1; d <= 3; d++){
        // Initialize LSQP
        lsqp_initialize( &data, &control, &status );
        // Set user-defined control options
        control.f_indexing = false; // C sparse matrix indexing
        // Start from 0
        real_wp_ x[] = {0.0, 0.0, 0.0};
        real_wp_ y[] = {0.0, 0.0};
        real_wp_ z[] = {0.0, 0.0, 0.0};
        switch(d){
            case 1: // sparse co-ordinate storage
```

```

        st = 'C';
        lsqp_import( &control, &data, &status, n, m,
                     "coordinate", A_ne, A_row, A_col, NULL );
        lsqp_solve_qp( &data, &status, n, m, w, x_0, g, f,
                       A_ne, A_val, c_l, c_u, x_l, x_u, x, c, y, z,
                       x_stat, c_stat );

        break;
    printf(" case %li break\n",d);
    case 2: // sparse by rows
        st = 'R';
        lsqp_import( &control, &data, &status, n, m,
                     "sparse_by_rows", A_ne, NULL, A_col, A_ptr );
        lsqp_solve_qp( &data, &status, n, m, w, x_0, g, f,
                       A_ne, A_val, c_l, c_u, x_l, x_u, x, c, y, z,
                       x_stat, c_stat );

        break;
    case 3: // dense
        st = 'D';
        int A_dense_ne = 6; // number of elements of A
        real_wp_ A_dense[] = {2.0, 1.0, 0.0, 0.0, 1.0, 1.0};
        lsqp_import( &control, &data, &status, n, m,
                     "dense", A_dense_ne, NULL, NULL, NULL );
        lsqp_solve_qp( &data, &status, n, m, w, x_0, g, f,
                       A_dense_ne, A_dense, c_l, c_u, x_l, x_u,
                       x, c, y, z, x_stat, c_stat );

        break;
    }
    lsqp_information( &data, &inform, &status );
    if(inform.status == 0){
        printf("%c:%6i iterations. Optimal objective value = %5.2f status = %li\n",
              st, inform.iter, inform.obj, inform.status);
    }else{
        printf("%c: LSQP_solve exit status = %li\n", st, inform.status);
    }
    //printf("x: ");
    //for( int i = 0; i < n; i++) printf("%f ", x[i]);
    //printf("\n");
    //printf("gradient: ");
    //for( int i = 0; i < n; i++) printf("%f ", g[i]);
    //printf("\n");
    // Delete internal workspace
    lsqp_terminate( &data, &control, &inform );
}
}

```

## 4.2 lsqptf.c

This is the same example, but now fortran-style indexing is used.

```

/* lsqptf.c */
/* Full test for the LSQP C interface using Fortran sparse matrix indexing */
#include <stdio.h>
#include <math.h>
#include "galahad_precision.h"
#include "galahad_cfunctions.h"
#include "galahad_lsqp.h"
int main(void) {
    // Derived types
    void *data;
    struct lsqp_control_type control;
    struct lsqp_inform_type inform;
    // Set problem data
    int n = 3; // dimension
    int m = 2; // number of general constraints
    real_wp_ g[] = {0.0, 2.0, 0.0}; // linear term in the objective
    real_wp_ f = 1.0; // constant term in the objective
    int A_ne = 4; // Jacobian elements
    int A_row[] = {1, 1, 2, 2}; // row indices
    int A_col[] = {1, 2, 2, 3}; // column indices
    int A_ptr[] = {1, 3, 5}; // row pointers
    real_wp_ A_val[] = {2.0, 1.0, 1.0, 1.0 }; // values
    real_wp_ c_l[] = {1.0, 2.0}; // constraint lower bound
    real_wp_ c_u[] = {2.0, 2.0}; // constraint upper bound
    real_wp_ x_l[] = {-1.0, - INFINITY, - INFINITY}; // variable lower bound
    real_wp_ x_u[] = {1.0, INFINITY, 2.0}; // variable upper bound
    real_wp_ w[] = {1.0,1.0,1.0};
    real_wp_ x_0[] = {0.0,0.0,0.0};
    // Set output storage
    real_wp_ c[m]; // constraint values
    int x_stat[n]; // variable status
    int c_stat[m]; // constraint status
}

```



```

char st;
int status;
printf(" Fortran sparse matrix indexing\n\n");
printf(" basic tests of qp storage formats\n\n");
for( int d=1; d <= 3; d++){
    // Initialize LSQP
    lsqp_initialize( &data, &control, &status );
    // Set user-defined control options
    control.f_indexing = true; // Fortran sparse matrix indexing
    // Start from 0
    real_wp_ x[] = {0.0,0.0,0.0};
    real_wp_ y[] = {0.0,0.0};
    real_wp_ z[] = {0.0,0.0,0.0};
    switch(d){
        case 1: // sparse co-ordinate storage
            st = 'C';
            lsqp_import( &control, &data, &status, n, m,
                "coordinate", A_ne, A_row, A_col, NULL );
            lsqp_solve_qp( &data, &status, n, m, w, x_0, g, f,
                A_ne, A_val, c_l, c_u, x_l, x_u, x, c, y, z,
                x_stat, c_stat );
            break;
        case 2: // sparse by rows
            st = 'R';
            lsqp_import( &control, &data, &status, n, m,
                "sparse_by_rows", A_ne, NULL, A_col, A_ptr );
            lsqp_solve_qp( &data, &status, n, m, w, x_0, g, f,
                A_ne, A_val, c_l, c_u, x_l, x_u, x, c, y, z,
                x_stat, c_stat );
            break;
        case 3: // dense
            st = 'D';
            int A_dense_ne = 6; // number of elements of A
            real_wp_ A_dense[] = {2.0, 1.0, 0.0, 0.0, 1.0, 1.0};
            lsqp_import( &control, &data, &status, n, m,
                "dense", A_dense_ne, NULL, NULL, NULL );
            lsqp_solve_qp( &data, &status, n, m, w, x_0, g, f,
                A_dense_ne, A_dense, c_l, c_u, x_l, x_u,
                x, c, y, z, x_stat, c_stat );
            break;
    }
    lsqp_information( &data, &inform, &status );
    if(inform.status == 0){
        printf("%c:%6i iterations. Optimal objective value = %5.2f status = %li\n",
            st, inform.iter, inform.obj, inform.status);
    }else{
        printf("%c: LSQP_solve exit status = %li\n", st, inform.status);
    }
    //printf("x: ");
    //for( int i = 0; i < n; i++) printf("%f ", x[i]);
    //printf("\n");
    //printf("gradient: ");
    //for( int i = 0; i < n; i++) printf("%f ", g[i]);
    //printf("\n");
    // Delete internal workspace
    lsqp_terminate( &data, &control, &inform );
}
}

```

