GALAHAD

USER DOCUMENTATION
GALAHAD Optimization Library version 4.0

## 1 SUMMARY

This package uses a preconditioned, projected-gradient method to solve the convex bound-constrained quadratic programming problem

$$
\text { minimize } q(\mathbf{x})=\frac{1}{2} \mathbf{x}^{T} \mathbf{H} \mathbf{x}+\mathbf{g}^{T} \mathbf{x}+f
$$

subject to the simple bound constraints

$$
x_{j}^{l} \leq x_{j} \leq x_{j}^{u}, \quad j=1, \ldots, n
$$

where the $n$ by $n$ symmetric, positive-definite matrix $\mathbf{H}$, the vectors $\mathbf{g}, \mathbf{x}^{l}, \mathbf{x}^{u}$ and the scalar $f$ are given. Any of the constraint bounds $x_{j}^{l}$ and $x_{j}^{u}$ may be infinite. Full advantage is taken of any zero coefficients in the matrix $\mathbf{H}$; the matrix need not be provided as there are options to obtain matrix-vector products involving $\mathbf{H}$ either by reverse communication or from a user-provided subroutine.

ATTRIBUTES - Versions: GALAHAD_BQP_single, GALAHAD_BQP_double. Uses: GALAHAD_CPU_time, GALAHAD_SYMBOLS, GALAHAD_SPACE, GALAHAD_SBLS, GALAHAD_QPT, GALAHAD_SPECFILE. Date: November 2009. Origin: N. I. M. Gould, Rutherford Appleton Laboratory. Language: Fortran 95 + TR 15581 or Fortran 2003.

## 2 HOW TO USE THE PACKAGE

Access to the package requires a USE statement such as
Single precision version
USE GALAHAD_BQP_single
Double precision version
USE GALAHAD_BQP_double
If it is required to use both modules at the same time, the derived types SMT_type, QPT_problem_type, NLPT_userdata-
 BQP_initialize, BQP_solve, BQP_terminate, (Section 2.3) and BQP_read_specfile (Section 2.7) must be renamed on one of the USE statements.

### 2.1 Matrix storage formats

When it is explicitly available, the Hessian matrix $\mathbf{H}$ may be stored in a variety of input formats.

### 2.1.1 Dense storage format

Since $\mathbf{H}$ is symmetric, only the lower triangular part (that is the part $h_{i j}$ for $1 \leq j \leq i \leq n$ ) need be held. The lowertriangular part of $\mathbf{H}$ is stored as a compact dense matrix by rows, that is, the values of the entries that occur before or on the diagonal of each row in turn are stored in order within an appropriate real one-dimensional array. That is component $i *(i-1) / 2+j$ of the storage array $\mathrm{H} \circ \mathrm{val}$ will hold the value $h_{i j}$ (and, by symmetry, $h_{j i}$ ) for $1 \leq j \leq i \leq n$.

[^0]
### 2.1.2 Sparse co-ordinate storage format

Only the nonzero entries of the lower triangular part of $\mathbf{H}$ is stored. For the $l$-th entry of the lower triangular part of $\mathbf{H}$, its row index $i$, column index $j$ and value $h_{i j}$ are stored in the $l$-th components of the integer arrays $\mathrm{H} \% \mathrm{row}, \mathrm{H} \% \mathrm{COl}$ and real array $H \% \mathrm{val}$, respectively. The order is unimportant, but the total number of entries $H \%$ ne is also required.

### 2.1.3 Sparse row-wise storage format

Again only the nonzero nonzero entries of the lower triangular part of $\mathbf{H}$ is stored, but this time they are ordered so that those that occur before or on the diagonal in row $i$ appear directly before those in row $i+1$. For the $i$-th row of $\mathbf{H}$, the $i$-th component of a integer array $\mathrm{H} \circ \mathrm{ptr}$ holds the position of the first entry in this row, while $\mathrm{H} \circ \mathrm{ptr}(n+1)$ holds the total number of entries plus one. The column indices $j$ and values $h_{i j}$ of the entries occurring before or on the diagonal in the $i$-th row are stored in components $l=H \% p t r(i), \ldots, H \% p t r(i+1)-1$ of the integer array $\mathrm{H} \% \mathrm{col}$, and real array $H \% v a l$, respectively.

For sparse matrices, this scheme almost always requires less storage than its predecessor.

### 2.1.4 Diagonal storage format

If $\mathbf{H}$ is diagonal (i.e., $h_{i j}=0$ for all $1 \leq i \neq j \leq n$ ) only the diagonals entries $h_{i i}, 1 \leq i \leq n$, need be stored, and the first $n$ components of the array $\mathrm{H} \% \mathrm{val}$ may be used for the purpose.

### 2.2 The derived data types

Ten derived data types are accessible from the package.

### 2.2.1 The derived data type for holding matrices

The derived data type SMT_TYPE is used to hold the matrix $\mathbf{H}$. The components of SMT_TYPE used here are:
n is a scalar component of type default INTEGER, that holds the number of columns in the matrix.
ne is a scalar variable of type default INTEGER, that holds the number of matrix entries.
type is a rank-one allocatable array of type default CHARACTER, that is used to indicate the matrix storage scheme used. Its precise length and content depends on the type of matrix to be stored (see §2.2.2).
val is a rank-one allocatable array of type default REAL (double precision in GALAHAD_BQP_double) and dimension at least ne, that holds the values of the entries. Each pair of off-diagonal entries $h_{i j}=h_{j i}$ of a symmetric matrix $\mathbf{H}$ is represented as a single entry (see §2.1.1-2.1.3). Any duplicated entries that appear in the sparse co-ordinate or row-wise schemes will be summed.
row is a rank-one allocatable array of type default INTEGER, and dimension at least ne, that may hold the row indices of the entries. (see §2.1.2).
col is a rank-one allocatable array of type default INTEGER, and dimension at least ne, that may the column indices of the entries (see §2.1.2-2.1.3).
ptr is a rank-one allocatable array of type default INTEGER, and dimension at least $m+1$, that may hold the pointers to the first entry in each row (see §2.1.3).

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### 2.2.2 The derived data type for holding the problem

The derived data type QPT_problem_type is used to hold the problem. The components of QPT_problem_type are:
n is a scalar variable of type default INTEGER, that holds the number of optimization variables, $n$.
H is scalar variable of type SMT_TYPE that holds the Hessian matrix H. The following components are used:
Hotype is an allocatable array of rank one and type default CHARACTER, that is used to indicate the storage scheme used. If the dense storage scheme (see Section 2.1.1) is used, the first five components of $H \%$ type must contain the string DENSE. For the sparse co-ordinate scheme (see Section 2.1.2), the first ten components of H\%type must contain the string COORDINATE, for the sparse row-wise storage scheme (see Section 2.1.3), the first fourteen components of H\%type must contain the string SPARSE_BY_ROWS, and for the diagonal storage scheme (see Section 2.1.4), the first eight components of $H \% t y p e$ must contain the string DIAGONAL.

For convenience, the procedure SMT_put may be used to allocate sufficient space and insert the required keyword into H\%type. For example, if prob is of derived type BQP_problem_type and involves a Hessian we wish to store using the co-ordinate scheme, we may simply

CALL SMT_put( prob\%H\%type, 'COORDINATE', istat )
See the documentation for the GALAHAD package SMT for further details on the use of SMT_put.
H\%ne is a scalar variable of type default INTEGER, that holds the number of entries in the lower triangular part of $\mathbf{H}$ in the sparse co-ordinate storage scheme (see Section 2.1.2). It need not be set for any of the other three schemes.
$H \% v a l$ is a rank-one allocatable array of type default REAL (double precision in GALAHAD_BQP_double), that holds the values of the entries of the lower triangular part of the Hessian matrix $\mathbf{H}$ in any of the storage schemes discussed in Section 2.1.
$H \% r o w ~ i s ~ a ~ r a n k-o n e ~ a l l o c a t a b l e ~ a r r a y ~ o f ~ t y p e ~ d e f a u l t ~ I N T E G E R, ~ t h a t ~ h o l d s ~ t h e ~ r o w ~ i n d i c e s ~ o f ~ t h e ~ l o w e r ~ t r i a n g u-~$ lar part of $\mathbf{H}$ in the sparse co-ordinate storage scheme (see Section 2.1.2). It need not be allocated for any of the other three schemes.
$\mathrm{H} \% \mathrm{COl}$ is a rank-one allocatable array variable of type default INTEGER, that holds the column indices of the lower triangular part of $\mathbf{H}$ in either the sparse co-ordinate (see Section 2.1.2), or the sparse row-wise (see Section 2.1.3) storage scheme. It need not be allocated when the dense or diagonal storage schemes are used.
$H \% p t r$ is a rank-one allocatable array of dimension $n+1$ and type default INTEGER, that holds the starting position of each row of the lower triangular part of $\mathbf{H}$, as well as the total number of entries plus one, in the sparse row-wise storage scheme (see Section 2.1.3). It need not be allocated when the other schemes are used.

G is a rank-one allocatable array of dimension $n$ and type default REAL (double precision in GALAHAD_BQP_double), that holds the gradient $\mathbf{g}$ of the linear term of the quadratic objective function. The $j$-th component of G , $j=1, \ldots, n$, contains $\mathbf{g}_{j}$.
$\mathrm{f} \quad$ is a scalar variable of type default REAL (double precision in GALAHAD_BQP_double), that holds the constant term, $f$, in the objective function.

X_1 is a rank-one allocatable array of dimension $n$ and type default REAL (double precision in GALAHAD_BQP_double), that holds the vector of lower bounds $\mathbf{x}^{l}$ on the the variables. The $j$-th component of $\mathrm{X}_{-} 1, j=1, \ldots, n$, contains $\mathbf{x}_{j}^{l}$. Infinite bounds are allowed by setting the corresponding components of $\mathrm{X}_{-1}$ to any value smaller than -infinity, where infinity is a component of the control array control (see Section 2.2.3).

[^1]X_u is a rank-one allocatable array of dimension $n$ and type default REAL (double precision in GALAHAD_BQP_double), that holds the vector of upper bounds $\mathbf{x}^{u}$ on the variables. The $j$-th component of $\mathrm{X}_{-} u, j=1, \ldots, n$, contains $\mathbf{x}_{j}^{u}$. Infinite bounds are allowed by setting the corresponding components of X_u to any value larger than that infinity, where infinity is a component of the control array control (see Section 2.2.3).

X is a rank-one allocatable array of dimension n and type default REAL (double precision in GALAHAD_BQP_double), that holds the values $\mathbf{x}$ of the optimization variables. The $j$-th component of $\mathrm{X}, j=1, \ldots, n$, contains $x_{j}$.
z is a rank-one allocatable array of dimension $n$ and type default REAL (double precision in GALAHAD_BQP_double), that holds the values $\mathbf{z}$ of estimates of the dual variables corresponding to the simple bound constraints (see Section 4). The $j$-th component of $\mathrm{Z}, j=1, \ldots, n$, contains $z j$.

### 2.2.3 The derived data type for holding control parameters

The derived data type BQP _control_type is used to hold controlling data. Default values may be obtained by calling BQP_initialize (see Section 2.3.1), while components may also be changed by calling BQP_read_specfile (see Section 2.7.1). The components of BQP_control_type are:
error is a scalar variable of type default $\operatorname{TNTEGER}$, that holds the stream number for error messages. Printing of error messages in $B Q P \_$solve and $B Q P$ _terminate is suppressed if error $\leq 0$. The default is error $=6$.
out is a scalar variable of type default INTEGER, that holds the stream number for informational messages. Printing of informational messages in BQP_solve is suppressed if out $<0$. The default is out $=6$.
print_level is a scalar variable of type default INTEGER, that is used to control the amount of informational output which is required. No informational output will occur if print_level $\leq 0$. If print_level $=1$, a single line of output will be produced for each iteration of the process. If print_level $\geq 2$, this output will be increased to provide significant detail of each iteration. The default is print_level $=0$.
start_print is a scalar variable of type default INTEGER, that specifies the first iteration for which printing will be permitted in GALAHAD_BQP_solve. If start_print is negative, printing will be permitted from the outset. The default is start_print $=-1$.
stop_print is a scalar variable of type default INTEGER, that specifies the last iteration for which printing will be permitted in GALAHAD_BQP_solve. If stop_print is negative, printing will be permitted once it has been started by start_print. The default is stop_print $=-1$.
print_gap is a scalar variable of type default INTEGER. Once printing has been started, output will occur once every print_gap iterations. If print_gap is no larger than 1 , printing will be permitted on every iteration. The default is print_gap $=1$.
maxit is a scalar variable of type default INTEGER, that holds the maximum number of iterations which will be allowed in GALAHAD_BQP_solve. The default is maxit $=1000$.
cold_start is a scalar variable of type default INTEGER, that should be set to 0 if a warm start is required (with variables assigned according to B_stat, see below), and to any other value if the values given in prob\%X suffice. The default is cold_start $=1$.
ratio_cg_vs_sd is a scalar variable of type default INTEGER, that specifies the ratio of how many iterations use CG rather steepest descent will be attempted. The default is ratio_cg_vs_sd $=1$.
change_max is a scalar variable of type default INTEGER, that specifies the maximum number of per-iteration changes in the working set permitted when allowing subspace solution rather than steepest descent (see §4). The default is change_max $=2$.

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cg_maxit is a scalar variable of type default INTEGER, that holds the maximum number of conjugate-gradient iterations which will be allowed per main iteration in GALAHAD_BQP_solve. The default is cg_maxit $=1000$.
infinity is a scalar variable of type default REAL (double precision in GALAHAD_BQP_double), that is used to specify which constraint bounds are infinite. Any bound larger than infinity in modulus will be regarded as infinite. The default is infinity $=10^{19}$.
stop_p is a scalar variable of type default REAL (double precision in GALAHAD_BQP_double), that holds the required accuracy for the primal infeasibility (see Section 4). The default is stop_p $=u^{1 / 3}$, where $u$ is EPSILON ( 1.0 ) (EPSILON (1.0DO) in GALAHAD_BQP_double).
stop_d is a scalar variable of type default REAL (double precision in GALAHAD_BQP_double), that holds the required accuracy for the dual infeasibility (see Section 4). The default is stop_d $=u^{1 / 3}$, where $u$ is EPSILON (1.0) (EPSILON (1.0DO) in GALAHAD_BQP_double).
stop_c is a scalar variable of type default REAL (double precision in GALAHAD_BQP_double), that holds the required accuracy for the violation of complementarity slackness (see Section 4). The default is stop_c $=u^{1 / 3}$, where $u$ is EPSILON (1.0) (EPSILON(1.0D0) in GALAHAD_BQP_double).
identical_bounds_tol is a scalar variable of type default REAL (double precision in GALAHAD_BQP_double). Each pair of variable bounds $\left(x_{j}^{l}, x_{j}^{u}\right)$ that is closer than identical_bounds_tol will be reset to the average of their values, $\frac{1}{2}\left(x_{j}^{l}+x_{j}^{u}\right)$. The default is identical_bounds_tol $=u$, where $u$ is EPSILON (1.0) (EPSILON (1.0D0) in GALAHAD_BQP_double).
stop_cg_relative and stop_cg_absolute are scalar variables of type default REAL (double precision in GALAHAD_BQP_double), that hold the relative and absolute convergence tolerances for the conjugate-gradient iteration that occurs in the face of currently-active constraints when constructing the search direction. _stop_cg_relative $=0.01$ and stop_cg_absolute $=\sqrt{u}$, where $u$ is EPSILON(1.0) (EPSILON (1.0D0) in GALAHAD_BQP_double).
zero_curvature is a scalar variable of type default REAL (double precision in GALAHAD_BQP_double)that specifies the threshold below which any objective-function curvature encountered is regarded as zero. The default is zero_curvature $=10 u$, where $u$ is EPSILON (1.0) (EPSILON (1.0D0) in GALAHAD_BQP_double).
cpu_time_limit is a scalar variable of type default REAL (double precision in GALAHAD_BQP_double), that is used to specify the maximum permitted CPU time. Any negative value indicates no limit will be imposed. The default is cpu_time_limit = - 1.0 .
exact_arcsearch is a scalar variable of type default LOGICAL, that should be set. TRUE. if the exact generalized Cauchy point, the first estimate of the minimizer of the objective in the Cauchy direction within the feasible box, is required, and .FALSE. if an approximation suffices. The default is exact_arcsearch = .TRUE..
space_critical is a scalar variable of type default LOGICAL, that must be set .TRUE. if space is critical when allocating arrays and .FALSE. otherwise. The package may run faster if space_critical is .FALSE. but at the possible expense of a larger storage requirement. The default is space_critical $=$.FALSE..
deallocate_error_fatal is a scalar variable of type default LOGICAL, that must be set. TRUE. if the user wishes to terminate execution if a deallocation fails, and .FALSE. if an attempt to continue will be made. The default is deallocate_error_fatal = .FALSE..
prefix is a scalar variable of type default CHARACTER and length 30, that may be used to provide a user-selected character string to preface every line of printed output. Specifically, each line of output will be prefaced by the string prefix (2:LEN (TRIM (prefix) ) -1), thus ignoring the first and last non-null components of the supplied string. If the user does not want to preface lines by such a string, they may use the default prefix $=$ " ".

[^2]SBLS_control is a scalar variable of type SBLS_control_type whose components are used to control the factorization and/or preconditioner used, performed by the package GALAHAD_SBLS. See the documentation for GALAHAD_SBLS for further details.

### 2.2.4 The derived data type for holding timing information

The derived data type BQP_time_type is used to hold elapsed CPU times for the various parts of the calculation. The components of BQP_time_type are:
total is a scalar variable of type default REAL, that gives the total time spent in the package.
analyse is a scalar variable of type default REAL, that gives the time spent analysing the required matrices prior to factorization.
factorize is a scalar variable of type default REAL, that gives the time spent factorizing the required matrices.
solve is a scalar variable of type default REAL, that gives the time spent computing the search direction.

### 2.2.5 The derived data type for holding informational parameters

The derived data type $B Q P$ _inform_type is used to hold parameters that give information about the progress and needs of the algorithm. The components of $B Q P$ _inform_type are:
status is a scalar variable of type default INTEGER, that gives the exit status of the algorithm. See Section 2.6 for details.
alloc_status is a scalar variable of type default INTEGER, that gives the status of the last attempted array allocation or deallocation. This will be 0 if status $=0$.
bad_alloc is a scalar variable of type default CHARACTER and length 80, that gives the name of the last internal array for which there were allocation or deallocation errors. This will be the null string if status $=0$.
factorization_status is a scalar variable of type default INTEGER, that gives the return status from the matrix factorization.
iter is a scalar variable of type default INTEGER, that gives the number of iterations performed.
obj is a scalar variable of type default REAL (double precision in GALAHAD_BQP_double), that holds the value of the objective function at the best estimate of the solution found.
time is a scalar variable of type $B Q P$ _time_type whose components are used to hold elapsed CPU times for the various parts of the calculation (see Section 2.2.4).

SBLS_inform is a scalar variable of type SBLS_inform_type whose components provide information about the progress and needs of the factorization/preconditioner performed by the package GALAHAD_SBLS. See the documentation for GALAHAD_SBLS for further details.

### 2.2.6 The derived data type for holding problem data

The derived data type BQP_data_type is used to hold all the data for a particular problem, or sequences of problems with the same structure, between calls of $B Q P$ procedures. This data should be preserved, untouched, from the initial call to $B Q P$ _initialize to the final call to $B Q P$ _terminate.

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### 2.2.7 The derived data type for holding user data

The derived data type NLPT_userdata_type is available to allow the user to pass data to and from user-supplied matrixvector product and preconditioning subroutines (see Section 2.4). Components of variables of type NLPT_userdata_type may be allocated as necessary. The following components are available:
integer is a rank-one allocatable array of type default INTEGER.
real is a rank-one allocatable array of type default REAL (double precision in GALAHAD_BQP_double)
complex is a rank-one allocatable array of type default COMPLEX (double precision complex in GALAHAD_BQP_double). character is a rank-one allocatable array of type default CHARACTER.
logical is a rank-one allocatable array of type default LOGICAL.
integer_pointer is a rank-one pointer array of type default INTEGER.
real_pointer is a rank-one pointer array of type default REAL (double precision in GALAHAD_BQP_double)
complex_pointer is a rank-one pointer array of type default COMPLEX (double precision complex in GALAHAD_BQP_double).
character_pointer is a rank-one pointer array of type default CHARACTER.
logical_pointer is a rank-one pointer array of type default LOGICAL.

### 2.2.8 The derived data type for holding reverse-communication data

The derived data type $B Q P$ _reverse_type is used to hold data needed for reverse communication when this is required. The components of BQP_reverse_type are:
nz_v_start is a scalar variable of type default INTEGER, that may be used to hold the starting position in NZ_v (see below) of the list of indices of nonzero components of $\mathbf{v}$.
nz_v_end is a scalar variable of type default INTEGER, that may be used to hold the finishing position in NZ_v (see below) of the list of indices of nonzero components of $\mathbf{v}$.

NZ_v is a rank-one allocatable array of dimension $n$ and type default INTEGER, that may be used to hold the indices of the nonzero components of $\mathbf{v}$. If used, components NZ_v (nz_v_start: $\left.n z_{-} v \_e n d\right)$ of $V$ (see below) will be nonzero.

V is a rank-one allocatable array of dimension $n$ and type default REAL (double precision in GALAHAD_BQP_double), that is used to hold the components of the output vector $\mathbf{v}$.

PROD is a rank-one allocatable array of dimension $n$ and type default REAL (double precision in GALAHAD_BQP_double), that is used to record the components of the resulting vector $\mathbf{H v}$.
nz_prod_end is a scalar variable of type default INTEGER, that is used to record the finishing position in NZ_prod (see below) of the list of indices of nonzero components of $\mathbf{H v}$ if required.

NZ_prod is a rank-one allocatable array of dimension $n$ and type default INTEGER, that is used to record the list of indices of nonzero components of $\mathbf{H v}$ if required. Components NZ_prod (1:nz_prod_end) of PROD should then be nonzero.

[^3]
### 2.3 Argument lists and calling sequences

There are three procedures for user calls (see Section 2.7 for further features):

1. The subroutine $B Q P$ _initialize is used to set default values, and initialize private data, before solving one or more problems with the same sparsity and bound structure.
2. The subroutine $B Q P$ _solve is called to solve the problem.
 the private data, allocated by BQP_solve, at the end of the solution process.

We use square brackets [ ] to indicate OPTIONAL arguments.

### 2.3.1 The initialization subroutine

Default values are provided as follows:

```
CALL BQP_initialize( data, control )
```

data is a scalar INTENT (INOUT) argument of type BQP_data_type (see Section 2.2.6). It is used to hold data about the problem being solved.
control is a scalar INTENT (OUT) argument of type BQP_control_type (see Section 2.2.3). On exit, control contains default values for the components as described in Section 2.2.3. These values should only be changed after calling BQP_initialize.

### 2.3.2 The quadratic programming subroutine

The quadratic programming solution algorithm is called as follows:
CALL $B Q P \_$solve( prob, B_stat, data, control, inform, userdata[, reverse, eval_HPROD] )
prob is a scalar INTENT (INOUT) argument of type QPT_problem_type (see Section 2.2.2). It is used to hold data about the problem being solved. The user must allocate all the array components, and set values for all components, except perhaps prob\%H; if the effect of $\mathbf{H}$ is only available to form products via reverse communication (see reverse below) or with a user-supplied subroutine (see eval_HPROD below) prob\%H and it is not needed, but otherwise prob\%H should be set using whichever of the matrix formats described in Section 2.1 is appropriate for $\mathbf{H}$ for the user's application.

The components prob\% X and $\mathrm{prob} \%$ must be set to initial estimates of the primal variables, $\mathbf{x}$ and dual variables for the bound constraints, $\mathbf{z}$, respectively. Inappropriate initial values will be altered, so the user should not be overly concerned if suitable values are not apparent, and may be content with merely setting prob $\% \mathrm{X}=0.0$ and prob $\% Z=0.0$.

On exit, the components prob\% X and prob\% $Z$ will contain the best estimates of the primal variables $\mathbf{x}$, and dual variables for the bound constraints $\mathbf{z}$, respectively. Restrictions: prob\%n $>0$ and (if $\mathbf{H}$ is provided) prob\% $\% \%$ ne $\geq-2$. prob\%H_type $\in\{$ 'DENSE', 'COORDINATE', 'SPARSE_BY_ROWS', 'DIAGONAL' $\}$.

B_stat is a rank-one INTENT (INOUT) array argument of dimension prob\%n and type default INTEGER, that indicates which of the simple bound constraints are in the current working set. Possible values for B_stat ( $j$ ) , $j=1, \ldots$, prob\%n, and their meanings are
$<0$ the $j$-th simple bound constraint is in the working set, on its lower bound,
$>0$ the $j$-th simple bound constraint is in the working set, on its upper bound, and
0 the $j$-th simple bound constraint is not in the working set.

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Suitable values must be supplied if control $\%$ bqp_control $\%$ cold_start $=0$ on entry, but need not be provided for other input values of control\%cold_start. Inappropriate values will be ignored. On exit, B_stat will contain values appropriate for the ultimate working set.
data is a scalar INTENT (INOUT) argument of type BQP_data_type (see Section 2.2.6). It is used to hold data about the problem being solved. It must not have been altered by the user since the last call to BQP_initialize.
control is a scalar INTENT(IN) argument of type BQP_control_type (see Section 2.2.3). Default values may be assigned by calling $B Q P$ _initialize prior to the first call to $B Q P \_$solve.
inform is a scalar INTENT (INOUT) argument of type BQP_inform_type (see Section 2.2.5). On initial entry, the component status must be set to 1 , while other components need not be set. A successful call to $B Q P \_s o l v e ~ i s ~$ indicated when the component status has the value 0 . For other return values of status, see Sections 2.5 and 2.6 .
userdata is a scalar INTENT (INOUT) argument of type NLPT_userdata_type whose components may be used to communicate user-supplied data (see Section 2.2.7) to and from the OPTIONAL subroutine eval_HPROD (see below).
reverse is an OPTIONAL scalar INTENT (INOUT) argument of type BQP_reverse_type (see Section 2.2.8). It is used to communicate reverse-communication data between the subroutine and calling program. If reverse is PRESENT and eval\%HPROD (see below) is absent, the user should monitor inform\%status on exit (see Section 2.5).
eval_HPROD is an OPTIONAL user-supplied subroutine whose purpose is to evaluate the value of the product $\mathbf{H v}$ of the Hessian of the objective function with a given vector v. See Section 2.4.1 for details. If eval_HPROD is present, it must be declared EXTERNAL in the calling program. If eval_HPROD is absent, GALAHAD_BQP_solve will use reverse communication (see Section 2.5) to obtain Hessian-vector products if reverse is PRESENT or otherwise require that the user has provided all relevant components of prob\% H .

### 2.3.3 The termination subroutine

All previously allocated arrays are deallocated as follows:

```
CALL BQP_terminate( data, control, inform )
```

data is a scalar INTENT (INOUT) argument of type BQP_data_type exactly as for BQP_solve, which must not have been altered by the user since the last call to $B Q P$ _initialize. On exit, array components will have been deallocated.
control is a scalar INTENT (IN) argument of type BQP_control_type exactly as for BQP_solve.
inform is a scalar INTENT (OUT) argument of type BQP_inform_type exactly as for BQP_solve. Only the component status will be set on exit, and a successful call to BQP_terminate is indicated when this component status has the value 0 . For other return values of status, see Section 2.6.

### 2.4 Matrix-vector operations

### 2.4.1 Hessian-vector products via internal evaluation

If the argument eval_HPROD is present when calling GALAHAD_BQP_solve, the user is expected to provide a subroutine of that name to evaluate the product of the Hessian of the objective function $\mathbf{H}$ with a given vector $\mathbf{v}$. The routine must be specified as

```
SUBROUTINE eval_HPROD( status, userdata, V, PROD[, NZ_v, nz_v_start, nz_v_end, &
    NZ_prod, nz_prod_end] )
```

[^4]whose arguments are as follows:
status is a scalar INTENT (OUT) argument of type default INTEGER, that should be set to 0 if the routine has been able to evaluate the product $\mathbf{H v}$ and to a non-zero value if the evaluation has not been possible.
userdata is a scalar INTENT (INOUT) argument of type NLPT_userdata_type whose components may be used to communicate user-supplied data to and from the subroutine (see Section 2.2.7).

V is a rank-one INTENT (IN) array argument of type default REAL (double precision in GALAHAD_BQP_double) whose components contain the vector $\mathbf{v}$. If components nz_v_start, nz_v_end and NZ_v (see below) are PRESENT, only components NZ_v (nz_v_start:nz_v_end) of V will be nonzero and the remaining components of v should be ignored. Otherwise, all components of v should be presumed to be nonzero.

PROD is a rank-one INTENT (OUT) array argument of type default REAL (double precision in GALAHAD_BQP_double) whose components on output contain the required components of Hv. If components nz_prod_end and NZ_prod (see below) are PRESENT, only the nonzero components NZ_prod ( 1 : nz _prod_end) of PROD need be assigned. Otherwise, all components of PROD must be set.
nz_v_start is an OPTIONAL scalar variable of type default INTEGER, that, if PRESENT, holds the starting position in NZ_v of the list of indices of nonzero components of $\mathbf{v}$.
nz_v_end is an OPTIONAL scalar variable of type default INTEGER, that, if PRESENT, holds the finishing position in NZ_v of the list of indices of nonzero components of $\mathbf{v}$.

NZ_v is an OPTIONAL rank-one allocatable array of dimension $n$ and type default INTEGERthat, if PRESENT, holds the indices of the nonzero components of $\mathbf{v}$. If any of nz_v_start, nz_v_end and NZ_v are absent, all components of V are assumed to be nonzero.
nz_prod_end is an OPTIONAL scalar variable of type default INTEGER, that, if PRESENT, must be set to record the number of non-zeros in $\mathbf{H v}$.

NZ_prod is an OPTIONAL rank-one allocatable array of dimension $n$ and type default INTEGERthat, if PRESENT, must be set to record the list of indices of nonzero components of $\mathbf{H v}$. If either of nz_prod_end and NZ_prod are absent, all components of PROD should be set even if they are zero.

### 2.5 Reverse Communication Information

A positive value of inform\%status on exit from BQP_solve indicates that GALAHAD_BQP_solve is seeking further information-this will happen if the user has chosen to evaluate matrix-vector products by reverse communication. The user should compute the required information and re-enter GALAHAD_BQP_solve with inform\%status and all other arguments (except those specifically mentioned below) unchanged.

Possible values of inform\%status and the information required are
2. The user should compute the matrix-vector product $\mathbf{H v}$ using the vector $\mathbf{v}$ whose components are stored in reverse $\% \mathrm{~V}$ and store the required product in reverse $\%$ PROD.
3. The user should compute the matrix-vector product $\mathbf{H v}$ using the vector $\mathbf{v}$ whose nonzero components are stored in positions reverse\%NZ_v (reverse\%nz_v_start:reverse\%nz_v_end) of reverse\%v. The remaining components of reverse $\% \mathrm{~V}$ should be ignored.
4. The user should compute the nonzero components of the matrix-vector product $\mathbf{H v}$ using the vector $\mathbf{v}$ whose nonzero components are stored in positions reverse\%NZ_v(reverse\%nz_v_start:reverse\%nz_v_end) of reverse\%V. The remaining components of reverse\%V should be ignored. The nonzero components must occupy positions reverse\%NZ_prod (1:reverse\%nz_prod_end) of reverse\%PROD, and the components reverse\%NZ_prod and reverse\%nz_prod_end must be set. This return can only happen if control eexact_arcsearch is .TRUE..

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### 2.6 Warning and error messages

A negative value of inform\%status on exit from BQP_solve or BQP_terminate indicates that an error has occurred. No further calls should be made until the error has been corrected. Possible values are:
-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.
-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.
-3. One of the restrictions prob\%n $>0$ or the requirement that prob\%H_type contain its relevant string 'DENSE', 'COORDINATE', 'SPARSE_BY_ROWS' or 'DIAGONAL' when $\mathbf{H}$ is available, has been violated.
-4 . The bound constraints are inconsistent.
-9. The analysis phase of the factorization failed; the return status from the factorization package is given in the component inform\%factor_status.
-10. The factorization failed; the return status from the factorization package is given in the component inform\%factor_status.
-16. The problem is so ill-conditioned that further progress is impossible.
-17 . The step is too small to make further impact.
-18. Too many iterations have been performed. This may happen if control\%maxit is too small, but may also be symptomatic of a badly scaled problem.
-19. The CPU time limit has been reached. This may happen if control\%cpu_time_limit is too small, but may also be symptomatic of a badly scaled problem.
-20. The Hessian matrix $\mathbf{H}$ of the objective function appears to be indefinite.
-23. An entry from the strict upper triangle of $\mathbf{H}$ has been specified.

### 2.7 Further features

In this section, we describe an alternative means of setting control parameters, that is components of the variable control of type $B Q P$ _control_type (see Section 2.2.3), by reading an appropriate data specification file using the subroutine BQP_read_specfile. This facility is useful as it allows a user to change BQP control parameters without editing and recompiling programs that call BQP .

A specification file, or specfile, is a data file containing a number of "specification commands". Each command occurs on a separate line, and comprises a "keyword", which is a string (in a close-to-natural language) used to identify a control parameter, and an (optional) "value", which defines the value to be assigned to the given control parameter. All keywords and values are case insensitive, keywords may be preceded by one or more blanks but values must not contain blanks, and each value must be separated from its keyword by at least one blank. Values must not contain more than 30 characters, and each line of the specfile is limited to 80 characters, including the blanks separating keyword and value.

The portion of the specification file used by BQP_read_specfile must start with a "BEGIN BQP" command and end with an "END" command. The syntax of the specfile is thus defined as follows:

[^5]```
( .. lines ignored by BQP_read_specfile .. )
    BEGIN BQP
        keyword value
        ........
        keyword value
    END
( .. lines ignored by BQP_read_specfile .. )
```

where keyword and value are two strings separated by (at least) one blank. The "BEGIN BQP" and "END" delimiter command lines may contain additional (trailing) strings so long as such strings are separated by one or more blanks, so that lines such as

```
BEGIN BQP SPECIFICATION
```

and
END BQP SPECIFICATION
are acceptable. Furthermore, between the "BEGIN BQP" and "END" delimiters, specification commands may occur in any order. Blank lines and lines whose first non-blank character is ! or * are ignored. The content of a line after a ! or * character is also ignored (as is the ! or * character itself). This provides an easy manner to "comment out" some specification commands, or to comment specific values of certain control parameters.

The value of a control parameters may be of three different types, namely integer, logical or real. Integer and real values may be expressed in any relevant Fortran integer and floating-point formats (respectively). Permitted values for logical parameters are "ON", "TRUE", ". TRUE.", "T", "YES", "Y", or "OFF", "NO", "N", "FALSE", ". FALSE." and "F". Empty values are also allowed for logical control parameters, and are interpreted as "TRUE".

The specification file must be open for input when BQP_read_specfile is called, and the associated device number passed to the routine in device (see below). Note that the corresponding file is REWINDed, which makes it possible to combine the specifications for more than one program/routine. For the same reason, the file is not closed by BQP_read_specfile.

### 2.7.1 To read control parameters from a specification file

Control parameters may be read from a file as follows:

CALL BQP_read_specfile( control, device )
control is a scalar INTENT (INOUT) argument of type BQP_control_type (see Section 2.2.3). Default values should have already been set, perhaps by calling BQP_initialize. On exit, individual components of control may have been changed according to the commands found in the specfile. Specfile commands and the component (see Section 2.2.3) of control that each affects are given in Table 2.1.
device is a scalar INTENT (IN) argument of type default INTEGER, that must be set to the unit number on which the specfile has been opened. If device is not open, control will not be altered and execution will continue, but an error message will be printed on unit control\%error.

### 2.8 Information printed

If control\%print_level is positive, information about the progress of the algorithm will be printed on unit control$\%$ out. If control p print_level $=1$, a single line of output will be produced for each iteration of the process. For the initial-feasible-point phase, this will include values of the current primal and dual infeasibility, and violation of complementary slackness, the feasibility-phase objective value, the current steplength, the value of the barrier

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| command | componentof control | value type |
| :--- | :--- | :--- |
| error-printout-device | oerror | integer |
| printout-device | \%out | integer |
| print-level | \%print_level | integer |
| start-print | \%start_print | integer |
| stop-print | \%stop_print | integer |
| iterations-between-printing | \%print_gap | integer |
| maximum-number-of-iterations | \%maxit | integer |
| cold-start | \%cold_start | integer |
| ratio-of-cg-iterations-to-steepest-descent | \%ratio_cg_vs_sd | integer |
| max-change-to-working-set-for-subspace-solution | \%change_max | integer |
| maximum-number-of-cg-iterations-per-iteration | \%cg_maxit | integer |
| infinity-value | \%infinity | real |
| primal-accuracy-required | \%stop_p | real |
| dual-accuracy-required | \%stop_d | real |
| complementary-slackness-accuracy-required | \%stop_c | real |
| identical-bounds-tolerance | \%identical_bounds_tol | real |
| cg-relative-accuracy-required | \%stop_cg_relative | real |
| cg-absolute-accuracy-required | \%stop_cg_absolute | real |
| zero-curvature-threshold | \%zero_curvature | real |
| maximum-cpu-time-limit | \%cpu_time_limit | real |
| exact-arcsearch-used | \%exact_arcsearch | logical |
| space-critical | \%space_critical | logical |
| deallocate-error-fatal | \%deallocate_error_fatal | logical |
| output-line-prefix | \%prefix | character |

Table 2.1: Specfile commands and associated components of control.
parameter, the number of backtracks in the linesearch and the elapsed CPU time in seconds. Once a suitable feasible point has been found, the iteration is divided into major iterations, at which the barrier parameter is reduced, and minor iterations, and which the barrier function is approximately minimized for the current value of the barrier parameter. For the major iterations, the value of the barrier parameter, the required values of dual feasibility and violation of complementary slackness, and the current constraint infeasibility are reported. Each minor iteration of the optimality phase results in a line giving the current dual feasibility and violation of complementary slackness, the objective function value, the ratio of predicted to achieved reduction of the objective function, the trust-region radius, the number of backtracks in the linesearch, the number of conjugate-gradient iterations taken, and the elapsed CPU time in seconds.

If control\%print_level $\geq 2$ this output will be increased to provide significant detail of each iteration. This extra output includes residuals of the linear systems solved, and, for larger values of control\%print_level, values of the primal and dual variables and Lagrange multipliers.

## 3 GENERAL INFORMATION

Use of common: None.
Workspace: Provided automatically by the module.
Other routines called directly: None.

[^6]Other modules used directly: $B Q P \_$solve calls the GALAHAD packages GALAHAD_CPU_time, GALAHAD_SYMBOLS, GALAHAD_SPACE, GALAHAD_SBLS, GALAHAD_QPT and GALAHAD_SPECFILE.

Input/output: Output is under control of the arguments control\%error, control\%out and control\%print_level.
Restrictions: prob\%n $>0$, prob\%H_type $\in\{$ 'DENSE', 'COORDINATE', 'SPARSE_BY_ROWS', 'DIAGONAL' $\}$ (if $\mathbf{H}$ is explicit).

Portability: ISO Fortran $95+$ TR 15581 or Fortran 2003. The package is thread-safe.

## 4 METHOD

The required solution $\mathbf{x}$ necessarily satisfies the primal optimality conditions

$$
\begin{equation*}
\mathbf{x}^{l} \leq \mathbf{x} \leq \mathbf{x}^{u} \tag{4.1}
\end{equation*}
$$

the dual optimality conditions

$$
\begin{equation*}
\mathbf{H x}+\mathbf{g}=\mathbf{z} \text { and } \mathbf{z}=\mathbf{z}^{l}+\mathbf{z}^{u} \tag{4.2}
\end{equation*}
$$

and

$$
\begin{equation*}
\mathbf{z}^{l} \geq 0 \text { and } \mathbf{z}^{u} \leq 0 \tag{4.3}
\end{equation*}
$$

and the complementary slackness conditions

$$
\begin{equation*}
\left(\mathbf{x}-\mathbf{x}^{l}\right)^{T} \mathbf{z}^{l}=0 \text { and }\left(\mathbf{x}-\mathbf{x}^{u}\right)^{T} \mathbf{z}^{u}=0, \tag{4.4}
\end{equation*}
$$

where the components of the vector $\mathbf{z}$ are known as the dual variables for the bounds, and where the vector inequalities hold componentwise. Projected-gradient methods iterate towards a point that satisfies these conditions by ultimately aiming to satisfy (4.2), while ensuring that (4.1), and (4.3) and (4.4) are satisfied at each stage. Appropriate norms of the amounts by which (4.1), (4.2) and (4.4) fail to be satisfied are known as the primal and dual infeasibility, and the violation of complementary slackness, respectively.

The method is iterative. Each iteration proceeds in two stages. Firstly, the so-called generalized Cauchy point for the quadratic objective is found. (The purpose of this point is to ensure that the algorithm converges and that the set of bounds which are satisfied as equations at the solution is rapidly identified.) Thereafter an improvement to the objective is sought using either a direct-matrix or truncated conjugate-gradient algorithm.

## References:

This is a specialised version of the method presented in
A. R. Conn, N. I. M. Gould and Ph. L. Toint (1988). Global convergence of a class of trust region algorithms for optimization with simple bounds. SIAM Journal on Numerical Analysis 25 433-460,

## 5 EXAMPLE OF USE

Suppose we wish to minimize $\frac{1}{2} x_{1}^{2}+x_{2}^{2}+x_{1} x_{2}+\frac{3}{2} x_{3}^{2}+2 x_{2}+1$ subject to the simple bounds $-1 \leq x_{1} \leq 1$ and $x_{3} \leq 2$. Then, on writing the data for this problem as

$$
\mathbf{H}=\left(\begin{array}{lll}
1 & 1 & \\
1 & 2 & \\
& & 3
\end{array}\right), \mathbf{g}=\left(\begin{array}{c}
0 \\
2 \\
0
\end{array}\right), \mathbf{x}^{l}=\left(\begin{array}{c}
-1 \\
-\infty \\
-\infty
\end{array}\right) \text { and } \mathbf{x}^{u}=\left(\begin{array}{c}
1 \\
\infty \\
2
\end{array}\right)
$$

in sparse co-ordinate format, we may use the following code:

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```
! THIS VERSION: GALAHAD 3.3 - 03/06/2021 AT 08:15 GMT.
    PROGRAM GALAHAD_BQP_EXAMPLE
    USE GALAHAD_BQP_double ! double precision version
    IMPLICIT NONE
    INTEGER, PARAMETER :: wp = KIND( 1.OD+0 ) ! set precision
    REAL ( KIND = wp ), PARAMETER :: infinity = 10.0_wp ** 20
    TYPE ( QPT_problem_type ) :: p
    TYPE ( BQP_reverse_type ) :: reverse
    TYPE ( BQP_data_type ) :: data
    TYPE ( BQP_control_type ) :: control
    TYPE ( BQP_inform_type ) :: inform
    TYPE ( GALAHAD_userdata_type ) :: userdata
    INTEGER :: S
    REAL ( KIND = wp ) :: t1, t2
    INTEGER, PARAMETER :: n = 3, h_ne = 4
    INTEGER, ALLOCATABLE, DIMENSION( : ) :: B_stat
! start problem data
    ALLOCATE( p%G( n ), p%X_l( n ), p%X_u( n ) )
    ALLOCATE( p%X( n ), p%Z( n ) )
    ALLOCATE( B_stat( n ) )
    p%new_problem_structure = .TRUE. ! new structure
    p%n = n ; p%f = 1.0_wp ! dimensions & objective constant
    p%G = (/ 0.0_wp, 2.0_wp, 1.0_wp /) ! objective gradient
    p%X_l = (/ - 1.0_wp, - infinity, 0.0_wp /) ! variable lower bound
    p%X_u = (/ infinity, 1.0_wp, 2.0_wp /) ! variable upper bound
    p%X = 0.0_wp ; p%Z = 0.0_wp ! start from zero
! sparse co-ordinate storage format
    CALL SMT_put( p%H%type, 'COORDINATE', s ) ! Co-ordinate storage for H
    ALLOCATE( p%H%val( h_ne ), p%H%row( h_ne ), p%H%col( h_ne ) )
    p%H%val = (/ 1.0_wp, 2.0_wp, 1.0_wp, 3.0_wp /) ! Hessian H
    p%H%row = (/ 1, 2, 2, 3 /) ! NB lower triangle
    p%H%col = (/ 1, 2, 1, 3 /) ; p%H%ne = h_ne
! problem data complete
    CALL BQP_initialize( data, control, inform ) ! Initialize control parameters
    control%infinity = infinity ! Set infinity
! control%print_level = 1 ! print one line/iteration
    inform%status = 1
    CALL BQP_solve( p, B_stat, data, control, inform, userdata )
    IF ( inform%status == 0 ) THEN ! Successful return
            WRITE( 6, "(' BQP: ', IO, ' iterations ', /, &
            & ' Optimal objective value =',
    & ES12.4, /, ' Optimal solution = ', ( 5ES12.4 ) )" ) &
        inform%iter, inform%obj, p%X
    ELSE ! Error returns
            WRITE( 6, "( ' BQP_solve exit status = ', IO ) " ) inform%status
            WRITE( 6, * ) inform%alloc_status, inform%bad_alloc
    END IF
    CALL BQP_terminate( data, control, inform ) ! delete workspace
    DEALLOCATE( p%G, p%X, p%X_l, p%X_u, p%Z, B_stat )
    DEALLOCATE( p%H%val, p%H%row, p%H%col, p%H%type )
    END PROGRAM GALAHAD_BQP_EXAMPLE
```

This produces the following output:

```
BQP: 2 iterations
Optimal objective value = -1.0000E+00
```

```
Optimal solution = 2.0000E+00 -2.0000E +00 0.0000E+00
```

The same problem may be solved holding the data in a sparse row-wise storage format by replacing the lines

```
! sparse co-ordinate storage format
! problem data complete
```

by

```
! sparse row-wise storage format
```

    CALL SMT_put( p\%H\%type, 'SPARSE_BY_ROWS', s ) ! Specify sparse-by-row
    ALLOCATE ( p\%H\%val (h_ne ), p\%H\%col(h_ne ), p\%H\%ptr(n + 1 ) )
    ALLOCATE ( p\%A\%val ( a_ne ), p\%A\%col( a_ne ), p\%A\%ptr (m+1 ) )
    p\%H\%val = (/ 1.0_wp, 2.0_wp, 1.0_wp, 3.0_wp /) ! Hessian H
    \(\mathrm{p} \% \mathrm{H} \% \mathrm{col}=(/ 1,2,3,3 /) \quad\) ! NB lower triangular
    \(\mathrm{p} \% \mathrm{H} \% \mathrm{ptr}=(/ 1,2,3,5 /) \quad\) ! Set row pointers
    ! problem data complete
or using a dense storage format with the replacement lines

```
! dense storage format
    CALL SMT_put( p%H%type, 'DENSE', s ) ! Specify dense
    ALLOCATE( p%H%val( n * ( n + 1 ) / 2 ) )
    p%H%val = (/ 1.0_wp, 0.0_wp, 2.0_wp, 0.0_wp, 1.0_wp, 3.0_wp /) ! Hessian
! problem data complete
```

respectively.
If instead $\mathbf{H}$ had been the diagonal matrix

$$
\mathbf{H}=\left(\begin{array}{lll}
1 & & \\
& 2 & \\
& & 3
\end{array}\right)
$$

but the other data is as before, the diagonal storage scheme might be used for $\mathbf{H}$, and in this case we would instead

```
CALL SMT_put( prob%H%type, 'DIAGONAL', s ) ! Specify dense storage for H
ALLOCATE( p%H%val( n ) )
p%H%val = (/ 1.0_wp, 2.0_wp, 3.0_wp /) ! Hessian values
```

The same problem may be solved using reverse communication with the following code:

```
! THIS VERSION: GALAHAD 3.3 - 03/06/2021 AT 08:15 GMT.
    PROGRAM GALAHAD_BQP_SECOND_EXAMPLE
    USE GALAHAD_BQP_double ! double precision version
    IMPLICIT NONE
    INTEGER, PARAMETER :: wp = KIND( 1.OD+0 ) ! set precision
    REAL ( KIND = wp ), PARAMETER :: infinity = 10.0_wp ** 20
    TYPE ( QPT_problem_type ) :: p
    TYPE ( BQP_reverse_type ) :: reverse
    TYPE ( BQP_data_type ) :: data
    TYPE ( BQP_control_type ) :: control
    TYPE ( BQP_inform_type ) :: inform
    TYPE ( GALAHAD_userdata_type ) :: userdata
    INTEGER :: nflag, i, j, k, l
    REAL ( KIND = wp ) :: v_j
    INTEGER, PARAMETER :: n = 3, h_ne = 4, h_all = 5
```

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```
! INTEGER, PARAMETER :: n = 3, h_ne = 3, h_all = 3
    INTEGER, ALLOCATABLE, DIMENSION( : ) :: B_stat, FLAG, ROW, PTR
    REAL ( KIND = wp ), ALLOCATABLE, DIMENSION( : ) :: VAL
! start problem data
    ALLOCATE( p%G( n ), p%X_l( n ), p%X_u( n ) )
    ALLOCATE( p%X( n ), p%Z( n ) )
    ALLOCATE( B_stat( n ), FLAG( n ) )
    ALLOCATE( VAL( h_all ), ROW( h_all ), PTR( n + 1 ) )
    p%new_problem_structure = .TRUE. ! new structure
    p%n = n ; p%f = 1.0_wp ! dimensions & objective constant
    p%G = (/ 0.0_wp, 2.0_wp, 1.0_wp /) ! objective gradient
    p%X_l = (/ - 1.0_wp, - infinity, 0.0_wp /) ! variable lower bound
    p%X_u = (/ infinity, 1.0_wp, 2.0_wp /) ! variable upper bound
    p%X = 0.0_wp ; p%Z = 0.0_wp ! start from zero
    PTR = (/ 1, 3, 5, 6 /) ! whole Hessian by rows
    ROW = (/ 1, 2, 1, 2, 3 /) ! for matrix-vector products
    VAL = (/ 1.0_wp, 1.0_wp, 1.0_wp, 2.0_wp, 3.0_wp /)
! problem data complete
    CALL BQP_initialize( data, control, inform ) ! Initialize control parameters
    control%infinity = infinity ! Set infinity
! control%print_level = 1 ! print one line/iteration
    control%maxit = 40 ! limit the # iterations
    control%print_gap = 100 ! print every 100 terations
! control%exact_gcp = .FALSE.
    nflag = 0 ; FLAG = 0
    inform%status = 1
1 0 ~ C O N T I N U E ~ ! ~ S o l v e ~ p r o b l e m ~ - ~ r e v e r s e ~ c o m m m u n i c a t i o n ~ l o o p
    CALL BQP_solve( p, B_stat, data, control, inform, userdata, reverse )
    SELECT CASE ( inform%status )
    CASE ( 0 ) ! Successful return
        WRITE( 6, "(' BQP: ', IO, ' iterations ', /, &
    & ' Optimal objective value =',
    & ES12.4, /, ' Optimal solution = ', ( 5ES12.4 ) )" ) &
        inform%iter, inform%obj, p%X
    CASE ( 2 ) ! compute H * v
        reverse%PROD = 0.0_wp
        DO j = 1, p%n
            v_j = reverse%V( j )
            DO k = PTR( j ), PTR( j + 1 ) - 1
                i = ROW( k )
                        reverse%PROD( i ) = reverse%PROD( i ) + VAL( k ) * v_j
            END DO
        END DO
        GO TO 10
    CASE ( 3 ) ! compute H * v for sparse v
        reverse%PROD = 0.0_wp
        DO l = reverse%nz_v_start, reverse%nz_v_end
            j = reverse%NZ_v( l ) ; v_j = reverse%V( j )
            DO k = PTR( j ), PTR( j + 1 ) - 1
                        i = ROW( k )
                        reverse%PROD( i ) = reverse%PROD( i ) + VAL( k ) * v_j
            END DO
        END DO
        GO TO 10
    CASE ( 4 ) ! compute H * v for very sparse v and record nonzeros
```

[^7]```
    nflag = nflag + 1
    reverse%nz_prod = 0
    DO l = reverse%nz_v_start, reverse%nz_v_end
    j = reverse%NZ_v( l ) ; v_j = reverse%V( j )
    DO k = PTR( j ), PTR( j + 1 ) - 1
        i = ROW( k )
        IF ( FLAG( i ) < nflag ) THEN
            FLAG( i ) = nflag
            reverse%PROD( i ) = VAL( k ) * v_j
            reverse%nz_prod_end = reverse%nz_prod_end + 1
            reverse%NZ_prod( reverse%nz_prod_end ) = i
        ELSE
            reverse%PROD( i ) = reverse%PROD( i ) + VAL( k ) * v_j
            END IF
        END DO
    END DO
    GO TO 10
CASE DEFAULT ! Error returns
    WRITE( 6, "( ' BQP_solve exit status = ', I6 ) " ) inform%status
END SELECT
CALL BQP_terminate( data, control, inform, reverse ) ! delete workspace
DEALLOCATE( p%G, p%X, p%X_l, p%X_u, p%Z, B_stat, FLAG, PTR, ROW, VAL )
END PROGRAM GALAHAD_BQP_SECOND_EXAMPLE
```

Notice that storage for the Hessian is now not needed. This produces the same output.
The same problem may also be solved by user-provided matrix-vector products as follows:

```
! THIS VERSION: GALAHAD 3.3 - 03/06/2021 AT 08:15 GMT.
    PROGRAM GALAHAD_BQP_THIRD_EXAMPLE
    USE GALAHAD_BQP_double ! double precision version
    IMPLICIT NONE
    INTEGER, PARAMETER :: wp = KIND( 1.OD+0 ) ! set precision
    REAL ( KIND = wp ), PARAMETER :: infinity = 10.0_wp ** 20
    TYPE ( QPT_problem_type ) :: p
    TYPE ( BQP_reverse_type ) :: reverse
    TYPE ( BQP_data_type ) :: data
    TYPE ( BQP_control_type ) :: control
    TYPE ( BQP_inform_type ) :: inform
    TYPE ( GALAHAD_userdata_type ) :: userdata
    INTEGER, PARAMETER :: n = 3, h_ne = 4, h_all = 5
    INTEGER, PARAMETER :: len_integer = 2 * n + 3 + h_all, len_real = h_all
    INTEGER, PARAMETER :: nflag = 2, st_flag = 2, st_ptr = st_flag + n
    INTEGER, PARAMETER :: st_row = st_ptr + n + 1, st_val = 0
    INTEGER, ALLOCATABLE, DIMENSION( : ) :: B_stat
    EXTERNAL :: HPROD
! start problem data
    ALLOCATE( p%G( n ), p%X_l( n ), p%X_u( n ) )
    ALLOCATE( p%X( n ), p%Z( n ) )
    ALLOCATE( userdata%integer( len_integer ), userdata%real( len_real ) )
    ALLOCATE( B_stat( n ) )
    p%new_problem_structure = .TRUE. ! new structure
    p%n = n ; p%f = 1.0_wp ! dimensions & objective constant
    p%G = (/ 0.0_wp, 2.0_wp, 1.0_wp /) ! objective gradient
    p%X_l = (/ - 1.0_wp, - infinity, 0.0_wp /) ! variable lower bound
    p%X_u = (/ infinity, 1.0_wp, 2.0_wp /) ! variable upper bound
    p%X = 0.0_wp ; p%Z = 0.0_wp ! start from zero
```

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```
! whole Hessian by rows for efficient matrix-vector products
    userdata%integer( st_ptr + 1 : st_ptr + n + 1 ) = (/ 1, 3, 5, 6 /)
    userdata%integer( st_row + 1 : st_row + h_all ) = (/ 1, 2, 1, 2, 3 /)
    userdata%real( st_val + 1 : st_val + h_all )
        = (/ 1.0_wp, 1.0_wp, 1.0_wp, 2.0_wp, 3.0_wp /)
! problem data complete
    CALL BQP_initialize( data, control, inform ) ! Initialize control parameters
    control%infinity = infinity ! Set infinity
! control%print_level = 1 ! print one line/iteration
    control%maxit = 40 ! limit the # iterations
    control%print_gap = 100 ! print every 100 terations
    control%exact_gcp = .FALSE.
    userdata%integer( 1 ) = n
    userdata%integer( nflag ) = 0
    userdata%integer( st_flag + 1 : st_flag + n ) = 0
    inform%status = 1
    CALL BQP_solve( p, B_stat, data, control, inform, userdata, &
                                    eval_HPROD = HPROD )
    IF ( inform%status == 0 ) THEN ! Successful return
        WRITE( 6, "( ' BQP: ', IO, ' iterations ', /,
        &
        & ' Optimal objective value =',
        & ES12.4, /, 'Optimal solution = ', ( 5ES12.4 ) )" ) &
        inform%iter, inform%obj, p%X
    ELSE
        WRITE( 6, "(' BQP_solve exit status = ', I6 ) " ) inform%status
    END IF
    CALL BQP_terminate( data, control, inform ) ! delete workspace
    DEALLOCATE( p%G, p%X, p%X_l, p%X_u, p%Z, B_stat )
    DEALLOCATE( userdata%integer, userdata%real )
    END PROGRAM GALAHAD_BQP_THIRD_EXAMPLE
        SUBROUTINE HPROD( status, userdata, V, PROD, NZ_v, nz_v_start, nz_v_end, &
                NZ_prod, nz_prod_end )
! compute the matrix-vector product H * v
        USE GALAHAD_USERDATA_double
        INTEGER, PARAMETER :: wp = KIND( 1.0D+0 )
        INTEGER, INTENT( OUT ) :: status
        TYPE ( GALAHAD_userdata_type ), INTENT( INOUT ) :: userdata
        REAL ( KIND = wp ), DIMENSION( : ), INTENT( IN ) :: V
        REAL ( KIND = wp ), DIMENSION( : ), INTENT( OUT ) :: PROD
        INTEGER, OPTIONAL, INTENT( IN ) :: nz_v_start, nz_v_end
        INTEGER, OPTIONAL, INTENT( INOUT ) :: nz_prod_end
        INTEGER, DIMENSION( : ), OPTIONAL, INTENT( INOUT ) :: NZ_v
        INTEGER, DIMENSION( : ), OPTIONAL, INTENT( INOUT ) :: NZ_prod
        INTEGER :: i, j, k, l, n, nflag, st_flag, st_ptr, st_row, st_val
        REAL ( KIND = wp ) :: v_j
        n = userdata%integer( 1 )
        nflag = 2
        st_flag = 2
        st_ptr = st_flag + n
        st_row = st_ptr + n + 1
        st_val = 0
! compute H * v for very sparse v and record nonzeros
    IF ( PRESENT( NZ_prod ) .AND. PRESENT( nz_prod_end ) ) THEN
        userdata%integer( nflag ) = userdata%integer( nflag ) + 1
```

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```
        nz_prod = 0
        DO l = nz_v_start, nz_v_end
            j = NZ_v( l ) ; v_j = V( j )
            DO k = userdata%integer( st_ptr + j ), &
                                    userdata%integer( st_ptr + j + 1 ) - 1
                i = userdata%integer( st_row + k )
                IF ( userdata%integer( st_flag + i ) < &
                                    userdata%integer( nflag ) ) THEN
                userdata%integer( st_flag + i ) = userdata%integer( nflag )
                PROD( i ) = userdata%real( st_val + k ) * v_j
                nz_prod_end = nz_prod_end + 1
                NZ_prod( nz_prod_end ) = i
            ELSE
                    PROD( i ) = PROD( i ) + userdata%real( st_val + k ) * v_j
            END IF
            END DO
        END DO
! compute H * v for sparse v
    ELSE IF ( PRESENT( NZ_v ) .AND. PRESENT( nz_v_start ) .AND. &
                PRESENT( nz_v_end ) ) THEN
            PROD = 0.0_wp
        DO l = nz_v_start, nz_v_end
            j = NZ_v( l ) ; v_j = V( j )
            DO k = userdata%integer( st_ptr + j ), &
                    userdata%integer( st_ptr + j + 1 ) - 1
            i = userdata%integer( st_row + k )
            PROD( i ) = PROD( i ) + userdata%real( st_val + k ) * v_j
            END DO
        END DO
! compute H * v
    ELSE
        PROD = 0.0_wp
        DO j = 1, n
            v_j = v( j )
            DO k = userdata%integer( st_ptr + j ), &
                                    userdata%integer( st_ptr + j + 1 ) - 1
                i = userdata%integer( st_row + k )
                PROD( i ) = PROD( i ) + userdata%real( st_val + k ) * v_j
            END DO
        END DO
    END IF
    status = 0
    END SUBROUTINE HPROD
```

Now notice how the matrix $\mathbf{H}$ is passed to the matrix-vector product evaluation routine via the integer and real components of the derived type userdata.

[^8]
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