

ACADOS / ACADO 2.0 / ???
Reference

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Chapter 1

Interfaces

1.1 OCP QP Interface

The interface describes an Optimal Control Problem (OCP) Quadratic Programming (QP) problem in the form

$$\min_{x,u} \sum_{n=0}^{N-1} \frac{1}{2} \begin{bmatrix} u_n \\ x_n \\ 1 \end{bmatrix}^T \begin{bmatrix} R_n & S_n & r_n \\ S_n^T & Q_n & q_n \\ r_n^T & q_n^T & 0 \end{bmatrix} \begin{bmatrix} u_n \\ x_n \\ 1 \end{bmatrix} + \frac{1}{2} \begin{bmatrix} x_N \\ 1 \end{bmatrix}^T \begin{bmatrix} Q_N & q_N \\ q_N^T & 0 \end{bmatrix} \begin{bmatrix} x_N \\ 1 \end{bmatrix} \quad (1.1)$$

$$s.t. \quad x_{n+1} = A_n x_n + B_n u_n + b_n \quad , \quad n = 0, \dots, N-1 \quad (1.2)$$

$$\underline{u}_n \leq u_n \leq \bar{u}_n \quad , \quad n = 0, \dots, N-1 \quad (1.3)$$

$$\underline{x}_n \leq x_n \leq \bar{x}_n \quad , \quad n = 0, \dots, N \quad (1.4)$$

$$\underline{d}_n \leq C_n x_n + D_n u_n \leq \bar{d}_n \quad , \quad n = 0, \dots, N-1 \quad (1.5)$$

$$\underline{d}_N \leq C_N x_N \leq \bar{d}_N \quad (1.6)$$

The C code interface looks like

```
int ocp_qp_SOLVERNAME(  
    int N, int *nx, int *nu, int *nb, int *ng,  
    double **A, double **B, double **b,  
    double **Q, double **S, double **R, double **q, double **r,  
    int **idxb, double **lb, double **ub,  
    double **C, double **D, double **ld, double **ud,  
    double **x, double **u,  
    struct ocp_qp_SOLVERNAME_args *args, double *work);
```

where SOLVERNAME is the name of the specific solver.

N [input] is the horizon length.

nx [input] is the vector of the state sizes n_u at the different stages, such that
nx[n] is the state size at stage n.

- nu** [input] is the vector of the input sizes n_x at the different stages, such that **nu**[**n**] is the input size at stage **n**.
- nb** [input] is the vector of the bound sizes n_b at the different stages, such that **nb**[**n**] is the bound size at stage **n**. The value of **nb**[**n**] is smaller or equal to **nx**[**n**]+**nu**[**n**].
- ng** [input] is the vector of the general polytopic constraint sizes n_g at the different stages, such that **ng**[**n**] is the general polytopic constraint size at stage **n**.
- A** [input] is the vector of size N of the pointers to the first element of the matrices A_n , such that **A**[**n**] is the pointer to the first element of the matrix A_n , and **A**[**n**][0] is the first element of the matrix A_n . The matrix referenced by the pointer **A**[**n**] is stored in column-major (or Fortran-like) order, in a vector of **nx**[**n**+1] \times **nx**[**n**] double-precision floating-point numbers.
- B** [input] is the vector of size N of the pointers to the first element of the matrices B_n , such that **B**[**n**] is the pointer to the first element of the matrix B_n , and **B**[**n**][0] is the first element of the matrix B_n . The matrix referenced by the pointer **B**[**n**] is stored in column-major (or Fortran-like) order, in a vector of **nx**[**n**+1] \times **nu**[**n**] double-precision floating-point numbers.
- b** [input] is the vector of size N of the pointers to the first element of the vectors b_n , such that **b**[**n**] is the pointer to the first element of the vector b_n , and **b**[**n**][0] is the first element of the vector b_n . The vector referenced by the pointer **b**[**n**] is stored in a vector of **nx**[**n**+1] \times 1 double-precision floating-point numbers.
- Q** [input] is the vector of size $N + 1$ of the pointers to the first element of the matrices Q_n , such that **Q**[**n**] is the pointer to the first element of the matrix Q_n , and **Q**[**n**][0] is the first element of the matrix Q_n . The matrix referenced by the pointer **Q**[**n**] is stored in column-major (or Fortran-like) order, in a vector of **nx**[**n**] \times **nx**[**n**] double-precision floating-point numbers.
- S** [input] is the vector of size N of the pointers to the first element of the matrices S_n , such that **S**[**n**] is the pointer to the first element of the matrix S_n , and **S**[**n**][0] is the first element of the matrix S_n . The matrix referenced by the pointer **S**[**n**] is stored in column-major (or Fortran-like) order, in a vector of **nu**[**n**] \times **nx**[**n**] double-precision floating-point numbers.
- R** [input] is the vector of size N of the pointers to the first element of the matrices R_n , such that **R**[**n**] is the pointer to the first element of the matrix R_n , and **R**[**n**][0] is the first element of the matrix R_n . The matrix

referenced by the pointer $\mathbf{R}[\mathbf{n}]$ is stored in column-major (or Fortran-like) order, in a vector of $\mathbf{nu}[\mathbf{n}] \times \mathbf{nu}[\mathbf{n}]$ double-precision floating-point numbers.

\mathbf{q} [input] is the vector of size $N + 1$ of the pointers to the first element of the vectors q_n , such that $\mathbf{q}[\mathbf{n}]$ is the pointer to the first element of the vector q_n , and $\mathbf{q}[\mathbf{n}][0]$ is the first element of the vector q_n . The vector referenced by the pointer $\mathbf{q}[\mathbf{n}]$ is stored in a vector of $\mathbf{nx}[\mathbf{n}] \times 1$ double-precision floating-point numbers.

\mathbf{r} [input] is the vector of size N of the pointers to the first element of the vectors r_n , such that $\mathbf{r}[\mathbf{n}]$ is the pointer to the first element of the vector r_n , and $\mathbf{r}[\mathbf{n}][0]$ is the first element of the vector r_n . The vector referenced by the pointer $\mathbf{r}[\mathbf{n}]$ is stored in a vector of $\mathbf{nu}[\mathbf{n}] \times 1$ double-precision floating-point numbers.

\mathbf{idxb} [input] is the vector of size $N + 1$ of the pointers to the first element of the integer vectors $idxb_n$ describing the indexes of the corresponding upper and lower bounds in \mathbf{lb} and \mathbf{ub} , such that $\mathbf{idxb}[\mathbf{n}]$ is the pointer to the index of the first bound at stage n , and $\mathbf{idxb}[\mathbf{n}][0]$ is index of the first bound at stage n . The indexes in $\mathbf{idxb}[\mathbf{n}]$ correspond to the position of the constrained components in the variables vector $\begin{bmatrix} u_n \\ x_n \end{bmatrix}$: therefore a bound on the first input component has index 0, a bound on the last input component has index $\mathbf{nu}[\mathbf{n}] - 1$, a bound on the first state component has index $\mathbf{nu}[\mathbf{n}]$ and a bound on the last state component has index $\mathbf{nu}[\mathbf{n}] + \mathbf{nx}[\mathbf{n}] - 1$. The vector referenced by the pointer $\mathbf{idxb}[\mathbf{n}]$ is stored in a vector of $\mathbf{nb}[\mathbf{n}] \times 1$ integer numbers.

\mathbf{lb} [input] is the vector of size $N + 1$ of the pointers to the first element of the vectors $\begin{bmatrix} u_n \\ x_n \end{bmatrix}$, such that $\mathbf{lb}[\mathbf{n}]$ is the pointer to the first element of the vector $\begin{bmatrix} u_n \\ x_n \end{bmatrix}$, and $\mathbf{lb}[\mathbf{n}][0]$ is the first element of the vector $\begin{bmatrix} u_n \\ x_n \end{bmatrix}$. The vector referenced by the pointer $\mathbf{lb}[\mathbf{n}]$ is stored in a vector of $\mathbf{nb}[\mathbf{n}] \times 1$ double-precision floating-point numbers.

\mathbf{ub} [input] is the vector of size $N + 1$ of the pointers to the first element of the vectors $\begin{bmatrix} \bar{u}_n \\ \bar{x}_n \end{bmatrix}$, such that $\mathbf{ub}[\mathbf{n}]$ is the pointer to the first element of the vector $\begin{bmatrix} \bar{u}_n \\ \bar{x}_n \end{bmatrix}$, and $\mathbf{ub}[\mathbf{n}][0]$ is the first element of the vector $\begin{bmatrix} \bar{u}_n \\ \bar{x}_n \end{bmatrix}$. The vector referenced by the pointer $\mathbf{ub}[\mathbf{n}]$ is stored in a vector of $\mathbf{nb}[\mathbf{n}] \times 1$ double-precision floating-point numbers.

\mathbf{C} [input] is the vector of size $N + 1$ of the pointers to the first element of the matrices C_n , such that $\mathbf{C}[\mathbf{n}]$ is the pointer to the first element of the matrix C_n , and $\mathbf{C}[\mathbf{n}][0]$ is the first element of the matrix C_n . The matrix

referenced by the pointer `C[n]` is stored in column-major (or Fortran-like) order, in a vector of `ng[n] × nx[n]` double-precision floating-point numbers.

D [input] is the vector of size N of the pointers to the first element of the matrices D_n , such that `D[n]` is the pointer to the first element of the matrix D_n , and `D[n][0]` is the first element of the matrix D_n . The matrix referenced by the pointer `D[n]` is stored in column-major (or Fortran-like) order, in a vector of `ng[n] × nu[n]` double-precision floating-point numbers.

ld [input] is the vector of size $N + 1$ of the pointers to the first element of the vectors \underline{d}_n , such that `ld[n]` is the pointer to the first element of the vector \underline{d}_n , and `ld[n][0]` is the first element of the vector \underline{d}_n . The vector referenced by the pointer `ld[n]` is stored in a vector of `ng[n] × 1` double-precision floating-point numbers.

ud [input] is the vector of size $N + 1$ of the pointers to the first element of the vectors \bar{d}_n , such that `ud[n]` is the pointer to the first element of the vector \bar{d}_n , and `ud[n][0]` is the first element of the vector \bar{d}_n . The vector referenced by the pointer `ud[n]` is stored in a vector of `ng[n] × 1` double-precision floating-point numbers.

x [output] is the vector of size $N + 1$ of the pointers to the first element of the vectors x_n , such that `x[n]` is the pointer to the first element of the vector x_n , and `x[n][0]` is the first element of the vector x_n . The vector referenced by the pointer `x[n]` is stored in a vector of `nx[n] × 1` double-precision floating-point numbers.

u [output] is the vector of size $N + 1$ of the pointers to the first element of the vectors u_n , such that `u[n]` is the pointer to the first element of the vector u_n , and `u[n][0]` is the first element of the vector u_n . The vector referenced by the pointer `u[n]` is stored in a vector of `nu[n] × 1` double-precision floating-point numbers.

args [input] is the pointer to a structure of type `ocp_qp_SOLVERNAME_args` that defines the arguments (as e.g. maximum number of iterations, minimum step size, ...) passed to the specific solver.

work [workspace] is the pointer to the working space used by the specific solver. The working space size (in doubles) is returned by a call to the function `ocp_qp_SOLVERNAME_workspace_double(int N, int *nx, int *nu, int *nb, int *ng, struct ocp_qp_solver_args *args)`.

Furthermore, the function returns an int, that is defined in the following enum (TODO change the names to something better!!!):

ACADOS_SUCCESS Solution successfully found.

ACADOS_MAXITER Maximum number of iterations reached.

ACADOS_MINSTEP Minimum step size reached (in IPs, probably unfeasible problem).

1.1.1 Examples

MPC problem

In the MPC problem, the initial state is fixed. This is modelled by choosing $\mathbf{nx}[0]=0$, i.e. not considering the initial state as an optimization variable. As a consequence, e.g. the matrix $\mathbf{A}[0]$ has size $\mathbf{nx}[1] \times 0$, the matrix $\mathbf{Q}[0]$ has size 0×0 , and the vector $\mathbf{q}[0]$ has size 0×1 . The information about the known value of x_0 and the matrix A_0 are used to compute the value of the vector $\mathbf{b}[0]$, that is initialized to $b_0 + A_0 \cdot x_0$.

1.2 Integrator Interface with Sensitivity Propagation

The C code interface looks like

```
int ocp_qp_SOLVERNAME(integrator_in *input,
                      struct integrator_SOLVERNAME_opts *opts, double *work);
```

where the input struct is defined as:

```
typedef struct integrator_in_{
    int nSteps, int nSystems, int nX, int nXA, int nP, int nOD,
    double *steps, int *nx, int *nxa, double **S0, double *mu,
    double **x, double **xa, double *p, double *od, double **xOut,
    double **sensOut, double *muOut, double *hessOut, int nOutputs,
    int *nGridOutputs, double **gridOutputs, double **muOutputs,
    double **valOutputs, double **sensOutputs, function_call *sys,
    function_call *out, double *variables} integrator_in;
```

where **SOLVERNAME** is the name of the specific solver. Additionally, a function struct has been defined as:

```
typedef struct function_call_{
    int dimIn, int dimOut, bool *sparsity, bool linear,
    void (*fun)(double*,double*), void (*jac)(double*,double*),
    bool forward, bool backward, bool hessian,
    void (*vde_forw)(double*,double*), void (*vde_adj)(double*,double*),
    void (*vde_hess)(double*,double*)} function_call;
```

If we like the latter embedding of structs, we can avoid some of the fields in the *integrator_in* struct by moving them somewhere else!

For fixed step integrators:

nSteps [input] is the number of integration steps (in case of no step size control).

steps [input] is the vector of size *nSteps*, containing the step size for each integration step.

Dimensions of the dynamic system(s):

nSystems [input] is the number of dynamic subsystems in order of dependency (states of subsystem *k* only depend on subsystems $1, 2, \dots, k - 1$ etc).

flag_linear [input] is the vector of size *nSystems*, containing the flag whether the subsystem is linear or not (0 or 1, should be part of options OR see the function_call struct).

nx [input] is the vector of size *nSystems*, containing the number of differential states for this integrator.

nxa [input] is the vector of size *nSystems*, containing the number of algebraic states for this integrator.

nX [input] is the **total** number of differential states for this integrator.

nXA [input] is the **total** number of algebraic states for this integrator.

nP [input] is the number of free parameters for this integrator (sensitivities with respect to these parameters are computed).

nOD [input] is the number of online data, i.e. fixed parameters for this integrator (no sensitivities are computed).

Sensitivity analysis for the integrator:

S0 [input] is the vector of size *nSystems* of the pointers to the matrices containing the forward seeds (the dimensions of these matrices are $nx[0] \times (nx[0] + np)$, $nx[1] \times (nx[0] + nx[1] + np)$, ... etc).

mu [input] is the vector of dimension *nX*, containing the backward seed.

flag_forward [input] is the flag for first order forward sensitivity propagation (0 or 1, should be part of options OR see the function_call struct).

flag_backward [input] is the flag for first order backward sensitivity propagation (0 or 1, should be part of options OR see the function_call struct).

flag_hessian [input] is the flag for second order Hessian propagation (0 or 1, should be part of options OR see the function_call struct).

Input and output data to the integrator:

x [input] is the vector of size *nSystems* of the pointers to the initial values of the differential states for this integrator.

xOut [output] is the vector of size $nSystems$ of the pointers to the end values of the differential states as a result of calling this integrator.

sensOut [output] is the vector of size $nSystems$ of the pointers to the end values of the forward sensitivities as a result of calling this integrator.

muOut [output] is the vector of size $nX + nP$ containing the backward derivatives as a result of calling this integrator.

hessOut [output] is the matrix of size $(nX + nP) \times (nX + nP)p$ containing the second order derivatives as a result of calling this integrator.

xa [input,output] is the vector of size $nSystems$ of the pointers to the initial **guess** of the algebraic states for this integrator (the guess can be updated by the integrator).

p [input] is the vector of size nP containing the free parameters for this integrator (sensitivities can be computed).

od [input] is the vector of size nOD containing the online data, i.e. fixed parameters for this integrator (no sensitivities are computed).

Extra outputs to be evaluated:

nOutputs [input] is the number of extra output functions (with each their own dimension).

dimOutputs [input] is the vector of size $nOutputs$, containing the dimension for each of the output functions (moved to function_call struct!).

sparsityOutputs [input] is the matrix of dimension $nOutputs \times (nX + nXA + nX)$, defining the sparsity pattern for each output function with respect to the differential states, the algebraic variables and the state derivatives (moved to function_call struct!).

nGridOutputs [input] is the vector of size $nOutputs$, containing the number of grid points on which each output function should be evaluated.

gridOutputs [input] is the vector of size $nOutputs$ of the pointers to the grid points on which each output function should be evaluated.

muOutputs [input] is the vector of size $nOutputs$ of the pointers to the backward seeds for each of the grid points defined by *gridOutputs*.

valOutputs [output] is the vector of size $nOutputs$ of the pointers to the outputs evaluated on the grid points defined by *gridOutputs*.

sensOutputs [output] is the vector of size $nOutputs$ of the pointers to the forward sensitivities of the outputs evaluated on the grid points defined by *gridOutputs*.

The functions (including AD) to be evaluated:

(This changed using the function_call struct!!)

rhs [input] the function pointer to evaluate the right-hand side.

jac [input] the function pointer to evaluate the full Jacobian.

vde_forw [input] the function pointer to evaluate the forward variational differential equations.

vde_adj [input] the function pointer to evaluate the adjoint equations.

vde_hess [input] the function pointer to evaluate the second order sensitivity equations.

out_vde_forw [input] is the vector of size *nOutputs*, containing the function pointers to evaluate the forward sensitivity equations for the outputs.

out_vde_adj [input] is the vector of size *nOutputs*, containing the function pointers to evaluate the adjoint equations for the output functions.

out_vde_hess [input] is the vector of size *nOutputs*, containing the function pointers to evaluate the second order sensitivity equations for the outputs.

Other memory:

variables [input,output] is the pointer to the warm variables space used by the specific solver from one call to the other. The variables space size (in doubles) is returned by a call to the function `integrator_SOLVERNAME_variables_double(integrator_in *input, struct integrator_solver_opts *opts)`.

opts [input] is the pointer to a structure of type `integrator_SOLVERNAME_opts` that defines the arguments (as e.g. maximum number of steps, minimum step size, desired tolerance, ...) passed to the specific solver.

work [workspace] is the pointer to the working space used by the specific solver. The working space size (in doubles) is returned by a call to the function `integrator_SOLVERNAME_workspace_double(integrator_in *input, struct integrator_solver_opts *opts)`.

Furthermore, the function returns an int, that is defined in the following enum (TODO change the names to something better!!!):

ACADOS_SUCCESS Solution successfully found.

ACADOS_MAXITER Maximum number of integration steps reached.

ACADOS_MINSTEP Minimum step size reached.