

The openMMF library: an open source software for multimode driven quantum systems

User manual

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Contents

1	How to build the library and compile the examples	2
2	What does the library calculate and how?	3
3	Use of the library	3
3.1	Declaration of the Hamiltonian	4
3.2	Hamiltonian components	5
3.3	Multimode Floquet matrix and diagonalisation	5
3.4	Time-evolution operator	5
3.5	Micromotion operator	6
3.6	Identifying the dressing modes	6
4	Typical performance	6
5	Default system types	7
5.1	Qubit	7
5.2	Spin of total angular momentum S_z	7
5.3	Ground state Alkali atoms.	8
6	List of examples	8
6.1	Python	8
6.2	C++	9
6.3	Fortran	9
7	C++ wrappers	10
8	Python wrappers	10
9	Bugs and known limitations	11
10	Acknowledgements	11
11	MODULES	12
11.1	Physical Constants	12
11.2	Arrays	13
11.3	Atomic properties	14
11.4	MKL	16
12	DERIVED TYPES (src/modes.f90)	17
13	COMPUTATIONAL SUBROUTINES	18
14	DRIVER SUBROUTINES	24
14.1	Utility subroutines	27

1 How to build the library and compile the examples

The openMMF source code includes a `Makefile` for compiling and building the library. The user must ensure that `make.inc` sets correctly the system's path to the LAPACK and (optionally) the MKL-intel libraries and linking/compiling options.

When the paths are set correctly, compiling the library requires invoking a single `make` command, which will build the library. To indicate whether the MKL library can be used, the user should indicate so in `make.inc`. The `make` building options are:

- `make` : Compiles the library and selected Fortran and C++ examples
- `make lib`: compiles the library including support for sparse matrices. This option requires the LAPACK and MKL-intel libraries.
- `make lib_lapack`: compiles the library without support for sparse matrices. This option requires LAPACK.
- `make Examples_lib`: compiles examples under the folder `examples/FORTRAN`, which uses LAPACK
- `make Examples_lib_sp`: compiles examples under the folder `examples/FORTRAN`, which uses the MKL library
- `make Examples_lib_c`: compiles examples under the folder `examples/CPP`, which uses LAPACK
- `make Examples_lib_c_sp`: compiles examples under the folder `examples/CPP`, which uses the MKL library
- `make` : run the options `lib` and `all_examples`.

All options of the `make` command produces static and dynamical libraries `libopenmmf.a` and `libopenmmf.so` in the folder `~/lib`. A number of `.mod` files produced with the building process are moved to the folder `~/include` and they are required for running the application.

The examples that only require the support of the LAPACK library are compiled with:

```
gfortran -o out SourceCode.f90 -I$(INCLUDE_OPENMMF) -L$(LIB_OPENMMF) -lopenmmf $(GFFLAGS)
g++      -o out SourceCode.cpp -I$(INCLUDE_OPENMMF) -L$(LIB_OPENMMF) -lopenmmf -lgfortran $(GFFLAGS)
```

while code that requires support from the MKL library is compiled by:

```
gfortran -o out SourceCode.f90 -I$(INCLUDE_OPENMMF) -L$(LIB_OPENMMF) -lopenmmf
                                           -L$(MKLLIBS) -I$(MKLINC) $(MKLFLAGS)
g++      -o out SourceCode.cpp -I$(INCLUDE_OPENMMF) -L$(LIB_OPENMMF) -lopenmmf -lgfortran
                                           -L$(MKLLIBS) -I$(MKLINC) $(MKLFLAGS)
```

The variables `MKLLIBS`, `MKLINC`, `GFFLAGS`, and `MKLFLAGS` are defined in the file `Makefile`. The environmental variables `INCLUDE_OPENMMF` and `LIB_OPENMMF` indicate the paths to the include and library directories of the openMMF library. Compilation of C++ source code follows the same formula, using the additional flag `-lgfortran` and the corresponding compiler `g++`. For an explicit example of usage in this case see the building script under `examples/CPP/Makefile`.

When running applications, the environmental variable `LD_LIBRARY_PATH` must indicate the path to such a library, which can be done with the shell command:

```
export LD_LIBRARY_PATH="/opt/intel/compilers_and_libraries/linux/mkl/lib/intel64:./lib"
```

which assumes default folder location of the MKL library and that `libopenmmf` is located in `./lib`.

2 What does the library calculate and how?

The library can be used to calculate the time-evolution operator, $U(t', t)$, $t' > t$, of systems whose Hamiltonian has the form:

$$H = \sum_{i,j}^D E_{i,j} |i\rangle \langle j| + \sum_{i,j}^D \sum_{\ell=1}^N \sum_{n \in \mathbb{Z}} V_{i,j}^{\ell,n} e^{in\omega_\ell t} |i\rangle \langle j| + \text{h.c.} \quad (1)$$

where D is the dimension of the Hilbert space, $E_{i,j}$ defines a static component of H , $V_{i,j}^{\ell,n}$ is the coupling between the states i and j oscillating at frequency $n\omega_\ell$ (i.e. the n -th harmonic of the ℓ -th fundamental frequency ω_ℓ) and N is the number of incommensurate frequencies.

To calculate the time-evolution operator we generalise the Rotating (or Resonant) Wave Approximation (RWA), taking into account the complex time dependence of Eq. (1). For this, we rephrase the problem in terms of building a time-dependent unitary transformation, $U_F(t)$ to a new basis $\{|\bar{i}\rangle\}$, that leads to a *time-independent* and diagonal Hamiltonian, \bar{H} . The operator $U_F(t)$ is called the *micromotion operator* and the new basis is a generalised multimode definition of the *dressed basis*. After applying the standard quantum-mechanical transformation rule to the Schrödinger equation [1, 2], this condition becomes:

$$U_F^\dagger(t) [H(t) - i\hbar\partial_t] U_F(t) = \sum_{\bar{i}} \bar{E}_{\bar{i}} |\bar{i}\rangle \langle \bar{i}| \quad (2)$$

where $\bar{E}_{\bar{i}}$ is the eigen-energy of the dressed state $|\bar{i}\rangle$, which is an eigenstate of the static Hamiltonian \bar{H} .

Importantly, in the basis of states defined by this transformation the time evolution operator is diagonal and has the form:

$$\bar{U}(t', t) = \sum_{\bar{i}} e^{-i\bar{E}_{\bar{i}}(t'-t)} |\bar{i}\rangle \langle \bar{i}| \quad (3)$$

which let us to calculate the time evolution operator in the original basis $\{|i\rangle\}$, just by inverting the transformation $U_F(t)$, according to [2]:

$$U(t', t) = U_F(t') \bar{U}(t', t) U_F(t) \quad (4)$$

To formulate a fully defined computational problem, we express the unitary transformation $U_F(t)$ as the multifrequency Fourier series [3]:

$$U_F(t) = \sum_{\vec{n}} U_{i,\bar{i}}^{\vec{n}} e^{-i\vec{\omega} \cdot \vec{n} t} |i\rangle \langle \bar{i}| \quad (5)$$

where $\vec{\omega} = (\omega_1, \omega_2, \dots, \omega_N)$ and \vec{n} is a N -dimensional vector of integers. After plugging this expansion in Eq. (2) and performing an integral over time, we obtain a fully defined eigenproblem for the eigenvalues $\bar{E}_{\bar{i}}$ and Fourier components of the unitary transformation $U_{i,\bar{i}}^{\vec{n}}$:

$$\sum_j (E_{i,j} - \hbar \vec{n} \cdot \vec{\omega}) U_{j,\bar{i}}^{\vec{n}} + \sum_j \sum_{\vec{m}} \left[V_{i,j}^{\vec{m}} U_{j,\bar{i}}^{\vec{n}+\vec{m}} + V_{j,i}^{\vec{m}*} U_{j,\bar{i}}^{\vec{m}-\vec{n}} \right] = \bar{E}_{\bar{i}} U_{i,\bar{i}}^{\vec{n}} \quad (6)$$

where $\vec{n}_{\ell,m} = \vec{n} + mP_\ell$ with $P_\ell = (0, \dots, 1, \dots, 0)$ the projector at the ℓ -th position. To obtain a finite matrix representation of this problem we truncate the sum over the number of modes of the Fourier expansion Eq. (5).

This formulation to calculate the time-evolution operator is equivalent to the multimode Floquet representation of the Hamiltonian that introduces an extended Hilbert space $|E_i, \vec{n}\rangle$ [3, 4]. However, the semiclassical description presented here makes emphasis in the physically accessible states.

3 Use of the library

Here we illustrate the use of the library's functionality considering a qubit driven by two harmonic forces. The Hamiltonian of this system has the form:

$$H = \hbar\omega_0 S_z + \hbar\Omega_1 \cos(\omega_1 t) S_x + \hbar\Omega_{2,x} \cos(\omega_2 t) S_x + \hbar\Omega_{2,y} \cos(\omega_2 t) S_z \quad (7)$$

The Fortran and C++ source codes to find the time-evolution operator are in the files:

- `examples/FORTRAN/main_DressedQubit.f90` .
- `examples/CPP/main_dressedqubit.cpp` .

3.1 Declaration of the Hamiltonian

First of all, we should declare the two derived types:

```
TYPE(ATOM)                                ID
TYPE(MODE),      DIMENSION(:),  ALLOCATABLE :: FIELDS
```

The variable `ID` contains information about the type of system, such as the number of levels and their energy spectrum (see the declaration of `TYPE(ATOM)` in Section 12). The derived type `FIELDS` stores information required to build the components of the Hamiltonian as well as the explicit matrix representation of the couplings.

In this concrete example, the components of `ID` are initialised by calling the subroutine:

```
CALL FLOQUETINIT(ID, 'qubit', INFO)
```

The second argument indicates the type of system (here 'qubit'), which, is sufficient to initialise the variable `ID` :

```
ID%id_system = 1
ID%D_BARE    = 2
```

where `ID%D_BARE` indicates the dimension of the Hilbert space.

Additional options of the function `FLOQUETINIT` are presented in section 5, which are useful for initialising some frequently used physical systems with default parameters. When dealing with a general quantum system (i.e. with an arbitrary energy spectrum), there is no need to call this function and the values of a variable of `TYPE(ATOM)` must be initialised explicitly.

In the source code, the next relevant instruction is the definition of an integer vector that provides information about the number of driving frequencies. The integer array `MODES_NUM` is allocated with size 3, indicating that the system will be driven by two fundamental frequencies (corresponding to $\ell = 1, 2$ in Eq. 1), since the first component is reserved to the static part of the Hamiltonian. The values of the elements of this array indicate the number of driving harmonics of each frequency, which here we set to 1 (making $n = 1$ in Eq. 1).

The total number of driving frequencies is equal to the sum of all elements of the array `MODES_NUM`. The user then should allocate sufficient memory space to store each one of the matrix representations of the couplings $V^{\ell,n}$. This is done with the sequence of instructions:

```
TOTAL_FREQUENCIES = SUM(MODES_NUM, 1)
ALLOCATE(FIELDS(TOTAL_FREQUENCIES))
DO m=1, TOTAL_FREQUENCIES
    ALLOCATE(FIELDS(m)%V(ID%D_BARE, ID%D_BARE))
END DO
```

By default the first element of the array of `FIELDS` is reserved for the static component of the Hamiltonian, which includes the spectrum of the static system as diagonal elements of the matrix `FIELDS(1)%V`. The next step then consist in defining all components of the Hamiltonian. For a quantum system with arbitrary energy spectrum, each one of the matrices `FIELDS(m)%V` must be declared explicitly.

When dealing with spin systems, as in the present example, each component of the Hamiltonian can be written as:

$$V^{\ell,n} = e^{\phi_x} X S_x + e^{\phi_y} Y S_y + e^{\phi_z} Z S_z \quad (8)$$

where S_i is the angular momentum operator. Therefore, we need only six parameters (three phases and three amplitudes) to define the coupling matrices.

These six parameters should be declared explicitly for each one of the the driving modes, along with the frequency (`omega`) and the corresponding number of modes to be included in the Fourier expansion of the evolution operator (`N_Floquet`). The values of these parameters are initialised as follows:

Source code	Parameter in Eq. (8)
<code>FIELDS(1)%X=0.0</code>	X
<code>FIELDS(1)%Y=0.0</code>	Y
<code>FIELDS(1)%Z=1.0</code>	Z
<code>FIELDS(1)%phi_x=0.0</code>	ϕ_x
<code>FIELDS(1)%phi_y=0.0</code>	ϕ_y
<code>FIELDS(1)%phi_z=0.0</code>	ϕ_z
<code>FIELDS(1)%omega=0.0</code>	ω
<code>FIELDS(1)%N_Floquet= 0</code>	N

where we have a direct correspondence with Eq. (8). This set of instructions is repeated for each one of the driving fields, as can be seen in the source code.

3.2 Hamiltonian components

The instructions detailed before let us to build the matrix representation of the terms in Eq. (8). This is done simply by calling the subroutine:

```
CALL SETHAMILTONIANCOMPONENTS(ID,size(modes_num,1),total_frequencies,MODES_NUM,FIELDS,INFO)
```

which results in storing the coupling $V^{\ell,n}$ in the set of matrices `FIELDS(r)%V`, with $r=1, \text{total_frequencies}$.

We remind the user that when the system of interest is not one of the default types defined in section 3.1, the user must define explicitly and in full all the coupling matrices. For an example of this situation, see the source file `example/FORTRAN/main_lattice.f90`.

3.3 Multimode Floquet matrix and diagonalisation

Once the components of the Hamiltonian are defined (i.e the complete set of matrices `FIELDS(r)%V` has been initialised), the multimode Hamiltonian can be calculated calling the function:

```
CALL MULTIMODEFLOQUETMATRIX(ID,size(modes_num,1),total_frequencies,MODES_NUM,FIELDS,INFO)
```

As a result of this call, the system stores the full multimode Floquet matrix on the left-hand side of Eq. (6) in the matrix `H_FLOQUET`. This matrix is defined in the module `ARRAYS` and can be accessed (and modified!) by all computational routines that include this module. The size of this matrix is calculated internally and stored in the parameter `h_floquet_size`, which is also a global variable.

The library includes a subroutine to evaluate a sparse representation of this matrix, which results after invoking:

```
CALL MULTIMODEFLOQUETMATRIX_SP(ID,SIZE(MODES_NUM,1),total_frequencies,  
                                MODES_NUM,FIELDS,VALUES,ROW_INDEX,COLUMN,INFO)
```

Setting `INFO = 0`, this instruction produces the representation of the Floquet matrix `H_FLOQUET` in the three array variation of the Compressed Sparse Row (CSR) storage format. With `INFO=6`, we get the matrix stored as three arrays of equal length, corresponding to values, rows and columns positions. The non-zero values of the matrix are stored in the complex array `VALUES`, and the information about their location is encoded in the integer arrays `COLUMN` and `ROW_INDEX`. The size of these three arrays are evaluated internally.

The library includes wrappers to diagonalisation subroutines from the Lapack and the MKL-intel (for the sparse CSR representation) libraries. These functions are called using:

Lapack :

```
CALL LAPACK_FULLEIGENVALUES(U_F,SIZE(H_FLOQUET,1),E_FLOQUET,INFO)
```

MKL:

```
CALL MKLSPARSE_FULLEIGENVALUES(D_MULTIFLOQUET,SIZE(VALUES,1),VALUES,ROW_INDEX,  
                                COLUMN,E_L,E_R,E_FLOQUET,U_F,INFO)
```

In both cases, the eigenvalues are stored in the array `E_FLOQUET` and the eigenvectors are stored as columns of the matrix `U_F`. Remember that these eigenvectors correspond to the coefficients of the multimode Fourier decomposition of the micromotion operator Eq. (5). Invoking the MKL subroutine requires two additional parameter `E_L` and `E_R`: the lower and upper bounds of the interval to be searched for eigenvalues, respectively. The user is responsible to set meaningful values of both parameters.

The MKL sparse solvers requires the use of the variation of the Compressed sparse row (CSR) representation of the matrix, which is produced by sorting the array of `ROW` position, in such a way that all non-zero values of a given row become stored consecutively. This sorting is done using a `QUICK_SORT` algorithm by Robert Renka (OAK Ridge Natl. Lab, US) accessible from the Alan Miller's Fortran Software website [5]. Alternatively, calling the function `MULTIMODEFLOQUETMATRIX_SP` with the variable `INFO=6`, the user obtains the extended Hamiltonian represented by the three arrays of `value,column,row`, which can be used with a different eigenvalue solver for sparse matrices.

3.4 Time-evolution operator

With the full spectrum of `H_FLOQUET`, the time evolution operator between `T1` and `T2`, corresponding to Eq. (4), can be evaluated calling the function:

```
CALL MULTIMODETIMEEVOLUTINOPERATOR(SIZE(U_F,1),SIZE(MODES_NUM,1),MODES_NUM,U_F,E_FLOQUET,  
                                    ID%D_BARE,FIELDS,T1,T2,U_AUX,INFO)
```

which produces the matrix representation of the operator in Eq. (4). The time evolution operator is stored in the complex matrix `U_AUX`, whose size is equal to the number of bare states `ID%D_BARE`.

3.5 Micromotion operator

The micromotion operator is the time-dependent unitary transformation between the bare basis and the basis of state where the Hamiltonian is time-independent. Since we know the Fourier decomposition of this transformation via the diagonalisation of H_{FLOQUET} , we can evaluate the instantaneous transformation, e.g. at time $T1$, using the subroutine:

```
CALL MULTIMODEMICROMOTION(ID,SIZE(U_F,1),NM,MODES_NUM,U_F,E_FLOQUET,ID%D_BARE,
                           FIELDS_,T1,U_F1,INFO)
```

which produces the matrix representation of the operator in Eq. (5). This micromotion operator is stored in the square matrix U_{F1} of size $ID%D_BARE$.

3.6 Identifying the dressing modes

In several applications is useful to define a dressed basis of states and the openMMF library includes functions to simplify the evaluation of the evolution operator in this basis. For this, first the user should identify the subset of driving fields that define the dressed states. This is done using an integer array with as many components as dressing fields. The elements of this array indicate the indices of the fields corresponding to the array `modes_num`. For example, if there is only one dressing field and it corresponds to the second component of the array `MODES_NUM`, then the array that indicates the dressing field, `DRESSINFIELDS_INDICES`, must be:

```
INTEGER, DIMENSION(2) :: DRESSINGFIELDS_INDICES
DRESSINGFIELDS_INDICES(1) = 1 ! THE STATIC COMPONENT
DRESSINGFIELDS_INDICES(2) = 2 ! THE FIRST DRIVING FIELD, WHICH DRESSES THE SYSTEM
```

With this, the Fourier decomposition of the micromotion operator defining the dressed basis in Eq. (5), can be obtained simply by calling the function:

```
CALL MICROMOTIONFOURIERDRESSED BASIS(ID,DRESSINGFIELDS_INDICES,MODES_NUM,
                                     FIELDS,U_FD,E_DRESSED,INFO)
```

The Fourier components $U_{i,\bar{i}}^{\vec{n}}$ are stored in the matrix U_{FD} and the spectrum of dressed energies, $\bar{E}_{\vec{i}}$, are stored in the array $E_{DRESSED}$. With these two elements, we can calculate the micromotion operator of the dressed basis using the subroutine:

```
CALL MICROMOTIONDRESSED BASIS(ID,MODES_NUM,DRESSINGFIELDS_INDICES,FIELDS,U_FD,
                              E_DRESSED,T1,U_FD_1,INFO)
```

The micromotion operator at $T1$ is then stored as the square matrix U_{FD_1} . This set of instructions let us to evaluate the time-evolution operator in the dressed basis using the sequence:

```
CALL MICROMOTIONDRESSED BASIS(ID,MODES_NUM,DRESSINGFIELDS_INDICES,FIELDS,U_FD,
                              E_DRESSED,T1,U_F1,INFO)
CALL MICROMOTIONDRESSED BASIS(ID,MODES_NUM,DRESSINGFIELDS_INDICES,FIELDS,U_FD,
                              E_DRESSED,T2,U_F2,INFO)
```

```
! ---- CALCULATE THE TIME-EVOLUTION OPERATOR IN THE DRESSED
! ---- BASIS USING THE PREVIOUSLY CALCULATED IN THE BARE BASIS
```

```
U_AUX = MATMUL(TRANPOSE(CONJG(U_F2)),MATMUL(U_AUX,U_F1))
```

4 Typical performance

The efficiency of the library is limited by the capability to solve the eigenvalue problem of the multimode Floquet Hamiltonian. Fig. ??(a) compares the execution times for a single diagonalisation of the Floquet Hamiltonian Bosonic and Fermionic Hubbard models, as functions of the size of the matrix (red and blue data/lines). The source codes can be found in `examples/FORTRAN/MANYBODYHUBBARD/` and the model Hamiltonian are:

$$H_{Bosons}(t) = (t + \Delta t \cos(\omega t + \phi)) \sum_{\langle i,j \rangle} b_i b_j^\dagger + \frac{u}{2} \sum_j n_j (n_j - 1)$$

$$H_{Fermions}(t) = (t + \Delta t \cos(\omega t + \phi)) \sum_{i,\sigma} (f_{i,\sigma} f_{i+1,\sigma}^\dagger + f_{i+1,\sigma}^\dagger f_{i,\sigma}) + u \sum_j n_{j,\uparrow} n_{j,\downarrow}$$

and we use the parameters $t = 1, u = 0.5, \phi = 0.0, \omega = 1.0, \Delta t = 0.2$. To vary the size of the Floquet Hamiltonian we increase the number of Floquet manifolds.

For these tests, the library performs two tasks: evaluates the Floquet Hamiltonian and use LAPACK to obtain the spectrum (only once for each matrix size). The Hamiltonian corresponds to a periodically driven Hubbard model (with near-neighbour tunnelling and on-site interaction) and the matrix dimension is given by the number of Floquet manifolds.

Figure ??(b) shows the execution time required to evaluate a single time-step of the time-evolution operator of driven qubit with a noisy background. The Hamiltonian of this system is:

$$H = \hbar\omega_0 S_z + \hbar\Omega S_x \cos(\omega t + \phi) + \sum_{\ell} \Delta_i \cos(\ell\omega/Nt + \phi_i) \quad (9)$$

with N the number of frequencies contributing to the noisy field, Δ_i and ϕ_i random amplitudes and phases, respectively. With this, the noise frequency is limited up to the driving frequency ω . The matrix size reflects the number of components of the noisy driving.

In this case we use a DRIVER subroutine and build a sparse representation of the multimode Hamiltonian. In contrast to the previous test, we observe that the matrix building and diagonalisation take about the same fraction of the total time. However, a direct comparison of the total times between these two examples is not straightforward since the structure of the Floquet matrix of the two problems is significantly different.

These results give us a rough idea of the scaling of the execution time with the matrix size, as well as the possible performance bottlenecks. These tests were performed using a DELL XPS with an Intel Core i7 vPro @ 2.7 GHz and 8GB of RAM. The machine runs on Linux Mint 17.2 Rafaela, with gcc-4.9.4

5 Default system types

The openMMF library defines three different system types by default. These are Qubit, Spin and the ground state of several alkali atoms. A general system with D energy levels can be defined using the “spin” type and setting by hand all coupling matrices.

5.1 Qubit

This type represents a two level system, where the couplings with the oscillating field are of the form

$$V^{\ell,n} = X S_x e^{\phi_{\ell,x}} + Y S_y e^{\phi_{\ell,y}} + Z S_z e^{\phi_{\ell,z}} \quad (10)$$

with $S_i, i \in x, y, z$ the set of spin 1/2 angular momentum operators with $\hbar := 1$.

The corresponding derived type is initialised with the instruction:

```
FLOQUET_INIT(ID, 'qubit', INFO)
```

5.2 Spin of total angular momentum S_z

This type represents a system with $2S_z + 1$ equally spaced energy levels, where the couplings with oscillating fields are of the form:

$$V^{\ell,n} = X S_x e^{\phi_{\ell,x}} + Y S_y e^{\phi_{\ell,y}} + Z S_z e^{\phi_{\ell,z}} \quad (11)$$

with $S_i, i \in x, y, z$ the set of angular momentum operators with total spin S_z .

The corresponding derived type is initialised with the instruction:

```
FLOQUET_INIT(ID, 'spin', 2*Sz, INFO)
```

where the third argument is a DOUBLEPRECISION variable equal to the double of the projection of the total angular momentum.

The matrix representation of the angular momentum operators s_i is build using [6]:

$$\langle S, m' | S_x | S, m \rangle = \frac{1}{2} \left[\sqrt{(S-m)(S+m+1)} \delta_{m,m'+1} + \sqrt{(S+m)(S-m+1)} \delta_{m,m'-1} \right] \quad (12)$$

$$\langle S, m' | S_y | S, m \rangle = \frac{i}{2} \left[\sqrt{(S-m)(S+m+1)} \delta_{m,m'+1} - \sqrt{(S+m)(S-m+1)} \delta_{m,m'-1} \right] \quad (13)$$

$$\langle S, m' | S_z | S, m \rangle = m \delta_{m,m'} \quad (14)$$

5.3 Ground state Alkali atoms.

The effective model of an alkali atom consist of an electron with zero orbital angular momenta interacting with a static nucleus. These two particles interact via their magnetic moments, which define two manifolds of states corresponding to the total angular momenta $F = I \pm 1/2$. The library can be used to study inter and intra manifold dynamics.

For example, to study the dynamics with focus on the manifold with total angular momentum $F = I - 1/2$ ('L', for lower), the corresponding ID can be obtained by invoking:

```
CALL FLOQUET_INIT(ID,SPECIE_NAME,'L',INFO)
```

and the Hamiltonian components are built assuming the form:

$$V^{\ell,n} = \frac{\mu_B g_{I-1/2} B_x}{A} F_x e^{\phi_{\ell,x}} + \frac{\mu_B g_{I-1/2} B_y}{A} F_y e^{\phi_{\ell,y}} + \frac{\mu_B g_{I-1/2} B_z}{A} F_z e^{\phi_{\ell,z}} \quad (15)$$

Similarly, to study the dynamic of the manifold with $F = I + 1/2$ ('U', for upper), the system is initialised using:

```
CALL FLOQUET_INIT(ID,SPECIE_NAME,'U',INFO)
```

which assumes couplings of the form

$$V^{\ell,n} = \frac{\mu_B g_{I+1/2} B_x}{A} F_x e^{\phi_{\ell,x}} + \frac{\mu_B g_{I+1/2} B_y}{A} F_y e^{\phi_{\ell,y}} + \frac{\mu_B g_{I+1/2} B_z}{A} F_z e^{\phi_{\ell,z}} \quad (16)$$

The atomic gyromagnetic factor is defined by:

$$g_F = g_J * (F * (F + 1) - I * (I + 1) + J * (J + 1)) / (2 * F * (F + 1)) + g_I * (F * (F + 1) + I * (I + 1) - J * (J + 1)) / (2 * F * (F + 1))! \quad (17)$$

with $F = I \pm 1/2$. The matrix representation in each case follows the same procedure as in the case of a spin system.

Finally, when both manifolds are of interest, we should use:

```
CALL FLOQUET_INIT(ID,SPECIE_NAME,'B',INFO)
```

which prepares the system to initialise couplings of the form:

$$V^{\ell,n} = \frac{\mu_B B_x}{A} (g_J J_x + g_I I_x) e^{\phi_{\ell,x}} + \frac{\mu_B B_y}{A} (g_J J_y + g_I I_y) e^{\phi_{\ell,y}} + \frac{\mu_B B_z}{A} (g_J J_z + g_I I_z) e^{\phi_{\ell,z}} \quad (18)$$

In this case, the angular momentum operators are represented in the uncoupled basis $|I, J, m_I, m_J\rangle$, corresponding to a tensor multiplication of the matrices resulting from Eqs. (12)-(14) representing I_i and J_i .

In all these cases, after invoking the initialisation routine `FLOQUET_INIT` and defining the parameters of the couplings using the derived data type `TYPE(MODES)::FIELDS`, the matrix representation of each component of the Hamiltonian can be obtained with a single call to the subroutine:

```
CALL SETHAMILTONIANCOMPONENTS(ID,size(modes_num,1),total_frequencies,MODES_NUM,FIELDS,INFO)
```

where `total_frequencies` is the total number of driving fields (including a static component). For a complete example see the source code at `examples/FORTRAN/main_87Rb.f90`.

6 List of examples

6.1 Python

Build the examples running `make` in the folder `examples/PYTHON`.

<code>main_87Rb.py</code>	Evaluates the time-evolution and spectrum of a driven atom of 87Rb.
<code>main_DressedQubit.py</code>	Evaluates the time evolution of a dressed qubit in the bare and dressed basis.
<code>main_floquetinit.py</code>	Illustrates the optional systems.
<code>main_qubit_driver.py</code>	Evaluates the tim-evolution operator of a driven qubit using a DRIVER subroutine.
<code>main_qubit.py</code>	Evaluates the time-evolution operator of a driven qubit.
<code>main_qubit.sp.py</code>	Evaluates the time-evolution operator of driven qubit using the sparse representation of the Hamiltonian.
<code>openmmf.py</code>	Defines the Python wrappers.

6.2 C++

Build the examples running **make** in the folder **examples/CPP** .

main_DressedQubit.cpp	Evaluates the time evolution of a dressed qubit in the bare and dressed basis.
main_DressedQubit.SP.cpp	Evaluates the time evolution of a dressed qubit in the bare and dressed basis using the sparse representation of the Hamiltonian.
main_qubit_driver.cpp	Evaluates the time-evolution operator of a driven qubit using a DRIVER subroutine.
main_qubit.cpp	Evaluates the time-evolution operator of a driven qubit.
main_qubit_sp.cpp	Evaluates the time-evolution operator of driven qubit using the sparse representation of the Hamiltonian.
MultimodeFloquet.h	Prototype of the C++ wrapper functions.

6.3 Fortran

Build the examples running **make** in the folder **examples/FORTRAN** .

main_qubit_Shirley.f90	Evaluates the time-evolution operator of a driven qubit with parameters as in the classical paper by Jon Shirley.
main_qubit.f90	Evaluates the time-evolution operator of a driven qubit.
main_qubit.SP.f90	Evaluates the time-evolution operator of driven qubit using the sparse representation of the Hamiltonian.
main_qubit_DRIVER.f90	Evaluates the time-evolution operator of a driven qubit using a DRIVER subroutine.
main_QuantumInterference.f90	Evaluates the time-evolution of a bichromatically driven qubit.
main_qubit_noise.SP.f90	Evaluates the time-evolution operator of a driven qubit in a noisy environment. Uses a sparse representation of the Hamiltonian.
main_halfspin_DRIVER.f90	Evaluates the time-evolution operator of a two-level system (defined with 'spin') using a DRIVER subroutine.
main_spin_DRIVER.f90	Evaluates the time-evolution operator of a L-level system (defined with 'spin') using a DRIVER subroutine.
main_lattice.f90	evaluates the spectrum and micromotion of a driven lattice with two sites in the unit cell.
main_threesiteslattice.f90	evaluates the spectrum and micromotion of a driven lattice with three sites in the unit cell.
main_qubitLattice.f90	Evaluates the time-evolution of a bichromatically driven qubit as in Eq. (20), arxiv 1612.02143.
main_qubitLattice.SP.f90	Evaluates the time-evolution of a bichromatically driven qubit as in Eq. (20), arxiv 1612.02143, using a sparse representation of the Hamiltonian.
main_qubitLattice_DRIVER.f90	Evaluates the time-evolution of a bichromatically driven qubit as in Eq. (20), arxiv 1612.02143, using a DRIVER subroutine.
main_DressedQubit.f90	Evaluates the time evolution of a dressed qubit in the bare and dressed basis.
main_DressedQubitV2.f90	Evaluates the time evolution of a dressed qubit in the bare and dressed basis.
main_DressedQubit.SP.f90	Evaluates the time evolution of a dressed qubit in the bare and dressed basis using the sparse representation of the Hamiltonian.
main_DressedQubit_DRIVER.f90	Evaluates the time evolution of a dressed qubit in the bare and dressed basis using a DRIVER subroutine.
main_87Rb_RFMW_DRIVER.f90	Evaluates the time-evolution operator of 87Rb dressed by a RF field and driven by a MW field, using the dressed basis.
DoublePeakSearch.f90	Evaluates the time-evolution operator of 87Rb dressed by a RF field and driven by a MW field, using the dressed basis.
Manybody/main_MBH_Bosons.f90	Evaluates the spectrum of the Bosonic Hubbard model with time-dependent tunnelling.
Manybody/main_MBH_Fermions.f90	Evaluates the spectrum of the Fermionic Hubbard model with time-dependent tunnelling.

7 C++ wrappers

The library includes C++ interfaces for each one of the subroutines defined. These interfaces are declared in files with the same name as the ones containing the Fortran declarations, but with the particle `_C` appended before the ending extension `.f90`. Similarly, wrapper subroutine are named using the corresponding Fortran names and appending the particle `_c` at the end of the name. For example, the file `MultimodeFloquetTE_C.f90` is paired with the file `MultimodeFloquet.f90` and defines the subroutine `MULTIMODETIMEEVOLUTIONOPERATOR_C`, which is used in c++ using `multimodetimeevolutionoperator_c_` followed by the declared list of arguments (see example at `examples/CPP/main_qubit.cpp`).

The prototype of all function enabled for C++ are declared in the header file `include/MultimodeFloquet.h`. This scheme lets us to establish a line-by-line correspondence between the Fortran and C++ source codes.

8 Python wrappers

The library includes an interface to Python, which is implemented using CTYPES. Similarly to the C++ wrappers, for Python we define a function for each one of the functionalities of the library. This set of functions is declared in the file `src/openmmf.py`, which should be copied to the python working directory. The Python wrappers call the C++ wrappers which are compiled in the file `libmultimodefloquet.so`. The user should make sure that the correct path to this file is set in `src/openmmf.py`.

The python wrapper requires to define the equivalent of the `atom` and `mode` derived data types, which are:

```
class atom_c_T(ctypes.Structure):
    _fields_ = [
        ("id_system", c_int),
        ("d_bare", c_int)
    ]
```

and

```
class mode_c_T(ctypes.Structure):
    c_dcplx = ctypes.c_double*2
    _fields_ = [
        ("omega", c_double),
        ("x", c_dcplx),
        ("y", c_dcplx),
        ("z", c_dcplx),
        ("phi_x", c_double),
        ("phi_y", c_double),
        ("phi_z", c_double),
        ("N_Floquet", c_int)
    ]
```

respectively.

Each python wrapper function in `src/openmmf.py` takes numpy input parameters and define the set of ctype pointers required to call the corresponding C++ wrapper function. Let's examine the instruction to build the multimode Floquet Hamiltonian:

```
def multimodefloquetmatrix(id,modes_num,fields,info):

    id_p          = ctypes.pointer(id)
    nm            = c_int(modes_num.size)
    total_frequencies = c_int(np.sum(modes_num))
    info          = c_int(info)
    modes_num_p   = modes_num.ctypes.data_as(POINTER(c_int))
    fields_p      = ctypes.pointer(fields)
    h_floquet_size =
    openmmfC.multimodefloquetmatrix_c_python_(id_p,ctypes.byref(nm),
        ctypes.byref(total_frequencies),
    modes_num_p,fields_p,ctypes.byref(info))

    return h_floquet_size
```

The input parameter `mode_num` is a numpy array of integers that, as before, defines the number of driving mode and harmonics. Also, the input parameter `info` is a numpy integer. `id` is an instance of the class `atom_c_T` and `fields` is an instance of the class `modes_c_T`. This function calls: `openmmfC.multimodefloquetmatrix_c_python_` passing all parameters by reference.

The ctypes wrappers rely on the equivalence between the data types in Fortran, C++ and Python, which should be observed rigorously following the Table 1

Fortran	C++	ctypes	numpy
integer	int	ctypes.c_int	numpy.int32
double precision	double	ctypes.c_double	numpy.double
complex*16	std::complex <double>	c_dcmplx = ctypes.c_double*2	np.complex

Table 1: Equivalence between the numerical data types in Fortran, C++, Ctypes and numpy.

9 Bugs and known limitations

If you find any bug please contact the developing team using the github hosting link <https://github.com/openMMF/>

10 Acknowledgements

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References

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11 MODULES

In this section we provide the header of each one of the subroutines of the library, including the argument declaration, to help the user to identify the type of variable expected by each function.

11.1 Physical Constants

The module `physical_constants` defines the default values of commonly used parameters defining the Hamiltonian of atomic systems. The user can modify these values accessing the file in `src/Modules.f90`.

```
MODULE physical_constants
  IMPLICIT NONE
  DOUBLE PRECISION, PARAMETER :: pi          = 4.0*ATAN(1.0)
  DOUBLE PRECISION, PARAMETER :: e           = 1.602176462D-19
  DOUBLE PRECISION, PARAMETER :: h_P         = 6.62606957D-34
  DOUBLE PRECISION, PARAMETER :: hbar        = h_P/(2.0*4.0*ATAN(1.0))
  DOUBLE PRECISION, PARAMETER :: mu_B        = 9.27400968D-24
  DOUBLE PRECISION, PARAMETER :: k_B         = 1.3806488D-23
  DOUBLE PRECISION, PARAMETER :: mu_cero     = 12.566370614D-7
  DOUBLE PRECISION, PARAMETER :: epsilon_cero = 8.854187817D-12
  DOUBLE PRECISION, PARAMETER :: amu         = 1.660538921D-27
  DOUBLE PRECISION, PARAMETER :: g_t         = 9.8D0
  DOUBLE PRECISION, PARAMETER :: SB_ct       = 5.6704D-8
  COMPLEX*16,      PARAMETER :: J_IMAG       = DCMLPX(0.0D0,1.0D0)
  DOUBLE PRECISION, PARAMETER :: speedoflight = 299792458.0D0
  DOUBLE PRECISION :: TOTAL_TIME
END MODULE physical_constants
```

11.2 Arrays

The module `ARRAYS` provides global definitions of matrices. When using the module, the user cannot define variables using any of the names declared in this module.

```
MODULE ARRAYS
  DOUBLE PRECISION, DIMENSION(:,:), ALLOCATABLE :: Identity,j_x,j_y,j_z,I_x,I_y,I_z
  COMPLEX*16,      DIMENSION(:,:), ALLOCATABLE :: H_IJ,HAMILTONIAN
  COMPLEX*16,      DIMENSION(:,:), ALLOCATABLE :: H_FLOQUET,H_FLOQUET_COPY
  COMPLEX*16,      DIMENSION(:,:), ALLOCATABLE :: U_ZEEMAN
END MODULE ARRAYS
```

To deallocate these arrays, the user can invoke the call:

```
CALL DEALLOCATEALL(ID)
```

where the variable `TYPE(ATOM) ID` defines the type of problem.

11.3 Atomic properties

The ATOMIC_PROPERTIES module defines the default physical parameters of common alkali atomic species used in the context of ultracold atomic physics.

```
MODULE ATOMIC_PROPERTIES
  USE physical_constants
  IMPLICIT NONE
  DOUBLE PRECISION :: L=0.0D0, S = 0.5D0
  DOUBLE PRECISION :: mass_at = 87*amu
  DOUBLE PRECISION :: I,g_I,g_J
  DOUBLE PRECISION :: J,F,gf,mf
  DOUBLE PRECISION :: gF_2,gF_1,G_F
  DOUBLE PRECISION :: A,a_s,alpha_E
  INTEGER           :: Fup,Fdown,Ftotal
  INTEGER           :: Total_states_LSI
  CHARACTER(LEN=7) :: ID_name

  !87Rb
  DOUBLE PRECISION :: I_87Rb = 1.5D0
  DOUBLE PRECISION :: J_87Rb = 0.5D0
  DOUBLE PRECISION :: gJ_87Rb = 2.0D0
  DOUBLE PRECISION :: gI_87Rb = -0.000995D0
  DOUBLE PRECISION :: A_87Rb = 2*pi*hbar*3.417341D9
  DOUBLE PRECISION :: a_s_87Rb = 5.77D-9
  DOUBLE PRECISION :: alpha_E_87Rb = 2*pi*hbar*0.0794*1D-4
  INTEGER           :: Fup_87Rb = 2
  INTEGER           :: Fdown_87Rb = 1
  CHARACTER(LEN=7) :: ID_name_87Rb = "87Rb"

  !6Li
  DOUBLE PRECISION :: I_6Li = 1.0D0
  DOUBLE PRECISION :: J_6Li = 0.5D0
  DOUBLE PRECISION :: gJ_6Li = 2.0D0
  DOUBLE PRECISION :: gI_6Li = -0.000995D0
  DOUBLE PRECISION :: A_6Li = 2*pi*hbar*152.137D6
  DOUBLE PRECISION :: a_s_6Li = 5.77D-9
  DOUBLE PRECISION :: alpha_E_6Li = 2*pi*hbar*0.0794*1D-4
  INTEGER           :: Fup_6Li = 1
  INTEGER           :: Fdown_6Li = 1
  CHARACTER(LEN=7) :: ID_name_6Li = "6Li"

  !qubit
  DOUBLE PRECISION :: I_qubit = 0.0D0
  DOUBLE PRECISION :: J_qubit = 0.0D0
  DOUBLE PRECISION :: gJ_qubit = 1.0D0
  DOUBLE PRECISION :: gI_qubit = 0.0D0
  DOUBLE PRECISION :: A_qubit = 1.0D0
  DOUBLE PRECISION :: a_s_qubit = 0.0D0
  DOUBLE PRECISION :: alpha_E_qubit = 0.0D0
  INTEGER           :: Fup_qubit = 1
  INTEGER           :: Fdown_qubit = 1
  CHARACTER(LEN=7) :: ID_name_qubit = "qubit"

  !spin
  DOUBLE PRECISION :: I_spin = 0.0D0
  DOUBLE PRECISION :: J_spin = 0.0D0
  DOUBLE PRECISION :: gJ_spin = 1.0D0
  DOUBLE PRECISION :: gI_spin = 0.0D0
  DOUBLE PRECISION :: A_spin = 1.0D0
  DOUBLE PRECISION :: a_s_spin = 0.0D0
  DOUBLE PRECISION :: alpha_E_spin = 0.0D0
```

```
INTEGER      :: Fup_spin    = 1
INTEGER      :: Fdown_spin  = 1
CHARACTER(LEN=7) :: ID_name_spin = "spin"

!lattice
CHARACTER      :: PERIODIC
CHARACTER(LEN=7) :: ID_name_lattice = "lattice"

END MODULE ATOMIC_PROPERTIES
```

11.4 MKL

```
MODULE FEAST
  integer      fpm(128)
  real*8       Emin,Emax
  real*8       epsout
  integer      loop
  integer      M0 ! initial guess
  integer      M1 ! total number of eigenvalues found
  integer      info_FEAST
  real*8,      DIMENSION(:),  ALLOCATABLE :: E, RES ! vector of eigenvalues
  complex*16, DIMENSION(:,:), ALLOCATABLE :: X      ! matrix with eigenvectors
END MODULE FEAST
```


12 DERIVED TYPES (src/modes.f90)

The derived type defined

```
MODULE TYPES
```

```
TYPE :: MODE
  DOUBLE PRECISION :: OMEGA
  COMPLEX*16        :: X,Y,Z
  DOUBLE PRECISION :: phi_x,phi_y,phi_z
  INTEGER           :: N_Floquet
  COMPLEX*16, DIMENSION(:,:), ALLOCATABLE :: V
  COMPLEX*16, DIMENSION(:),   ALLOCATABLE :: VALUES
  INTEGER,    DIMENSION(:),   ALLOCATABLE :: ROW,COLUMN
END TYPE MODE
```

```
TYPE :: ATOM
  INTEGER :: id_system
  INTEGER :: D_BARE
  DOUBLE PRECISION, DIMENSION(:), ALLOCATABLE :: E_BARE
END TYPE ATOM
```

```
TYPE :: HARMONIC_FACTORS
  COMPLEX*16,DIMENSION(:,:), ALLOCATABLE :: U,U_r,U_AVG
  INTEGER,    DIMENSION(:),   ALLOCATABLE :: n
END type HARMONIC_FACTORS
```

```
END MODULE TYPES
```

13 COMPUTATIONAL SUBROUTINES

```

MODULE FLOQUETINITINTERFACE
  INTERFACE FLOQUETINIT
    MODULE PROCEDURE FLOQUETINIT_QUBIT, FLOQUETINIT_SPIN, FLOQUETINIT_ALKALI
  END INTERFACE FLOQUETINIT
contains
  SUBROUTINE FLOQUETINIT_QUBIT(ID,atomicspecie,INFO)
    ! ATOMICSPECIE: 87Rb,6Li,Cs,41K,qubit,lattice, SPIN
    ! MANIFOLD : "U" UPPER HYPERFINE MANIFOLD, "L" LOWER HYPERFINE MANIFOLD, "B" BOTH
    ! JTOTAL : IF ATOMICSPECIE .EQ. SPIN THEN JTOTAL IS THE TOTAL ANGULAR MOMENTUM OF THE SPIN
    !           IF ATOMICSPECIE .EQ. LATTICE, THEN JTOTAL IS THE NUMBER OF SITES
    ! calculate the dimension of the Hilbert space
    ! initialize all the matrices required for a full Floquet calculations
    ! Calculate the nuclear, electron and total angular momentum operators

    TYPE(ATOM),          INTENT(INOUT) :: ID
    CHARACTER (LEN=*), INTENT(IN)      :: ATOMICSPECIE
    INTEGER,             INTENT(INOUT) :: INFO

    ...

  SUBROUTINE FLOQUETINIT_SPIN(ID,atomicspecie,jtotal,info)
    ! ATOMICSPECIE: 87Rb,6Li,Cs,41K,qubit,lattice, SPIN
    ! MANIFOLD : "U" UPPER HYPERFINE MANIFOLD, "L" LOWER HYPERFINE MANIFOLD, "B" BOTH
    ! JTOTAL : IF ATOMICSPECIE .EQ. SPIN THEN JTOTAL IS THE TOTAL ANGULAR MOMENTUM OF THE SPIN
    !           IF ATOMICSPECIE .EQ. LATTICE, THEN JTOTAL IS THE NUMBER OF SITES

    ! calculate the dimension of the Hilbert space
    ! initialize all the matrices required for a full Floquet calculations
    ! Calculate the nuclear, electron and total angular momentum operators

    IMPLICIT NONE
    TYPE(ATOM),          INTENT(INOUT) :: ID
    CHARACTER (LEN=*), INTENT(IN)      :: ATOMICSPECIE
    DOUBLE PRECISION,    intent(in)    :: jtotal
    INTEGER,             INTENT(INOUT) :: INFO

    ...

  SUBROUTINE FLOQUETINIT_ALKALI(ID,atomicspecie,manifold,info)
    ! ATOMICSPECIE: 87Rb,6Li,Cs,41K,qubit,lattice, SPIN
    ! MANIFOLD : "U" UPPER HYPERFINE MANIFOLD, "L" LOWER HYPERFINE MANIFOLD, "B" BOTH
    ! JTOTAL : IF ATOMICSPECIE .EQ. SPIN THEN JTOTAL IS THE TOTAL ANGULAR MOMENTUM OF THE SPIN
    !           IF ATOMICSPECIE .EQ. LATTICE, THEN JTOTAL IS THE NUMBER OF SITES

    ! calculate the dimension of the Hilbert space
    ! initialize all the matrices required for a full Floquet calculations
    ! Calculate the nuclear, electron and total angular momentum operators

    TYPE(ATOM),          INTENT(INOUT) :: ID
    CHARACTER (LEN=*), INTENT(IN)      :: ATOMICSPECIE
    CHARACTER (LEN=1), INTENT(IN)      :: MANIFOLD
    INTEGER,             INTENT(INOUT) :: INFO

    ...

END MODULE FLOQUETINITINTERFACE

```

```

SUBROUTINE SETHAMILTONIANCOMPONENTS(ID,NM,NF,MODES_NUM,FIELD,INFO)
! ID  tYPE OF ATOM
! MODES_NUM, VECTOR. THE SIZE OF THE VECTOR TELL US THE NUMBER OF
!           FREQUENCIES, AND THE VALUE OF EACH COMPONENT INDICATES
!           THE NUMBER OF HARMONICS OF EACH FREQUENCI
! FIELDS : IN AND OUTPUT THE MATRICES
! INFO

```

```

USE ARRAYS
USE ATOMIC_PROPERTIES
USE TYPES
USE SUBINTERFACE_LAPACK ! write_matrix interface

```

```

IMPLICIT NONE
INTEGER,          INTENT(IN)      :: NM,NF
TYPE(ATOM),       INTENT(IN)      :: ID
INTEGER,    DIMENSION(NM), INTENT(IN)  :: MODES_NUM
TYPE(MODE), DIMENSION(NF), INTENT(INOUT) :: FIELD
INTEGER,          INTENT(INOUT)  :: INFO

```

```

SUBROUTINE F_representation(Fx,Fy,Fz,Ftotal)

```

```

USE FUNCIONES

```

```

IMPLICIT NONE
DOUBLE PRECISION, DIMENSION(:,:), INTENT(OUT):: Fx,Fy,Fz
DOUBLE PRECISION, INTENT(IN) :: Ftotal
!INTEGER, INTENT(IN) :: Ftotal_

```

```

!DOUBLE PRECISION
INTEGER k,p,N_k
double precision k_!,Ftotal

```

```

Fx = 0.0
Fy = 0.0
Fz = 0.0

```

```

SUBROUTINE I_and_J_representations(j_x,j_y,j_z,I_x,I_y,I_z,L,S,I)

```

```

USE FUNCIONES

```

```

IMPLICIT NONE
DOUBLE PRECISION, DIMENSION(:,:),INTENT(INOUT) :: j_x,j_y,j_z,I_x,I_y,I_z
DOUBLE PRECISION, INTENT(IN) :: L,S,I

```

```

SUBROUTINE MULTIMODETIMEEVOLUTINOPERATOR(D,NM,MODES_NUM,U_F_MODES,E_MULTIFLOQUET,
D_BARE,FIELD,T1,T2,U,INFO)

```

```

! TIME EVOLUTION OPERATOR OF A MULTIMODE DRESSED SYSTEM.

```

```

! THE EVOLUTION OPERATOR IS WRITTEN IN THE BASIS USED TO EXPRESS THE
! MULTIMODE FLOQUET HAMILTONIAN
! U : MATRIX OF AMPLITUDES OF PROBABILITIES FOR TRANSITIONS BETWEEN T1 TO T2
!!$ D (IN) : DIMENSION OF THE EXTENDED HILBERT SPACE
!!$ (SIZE OF THE MULTIMODE FLOQUET MATRIX)
!!$ NM (IN) : NUMBER OF MODES
!!$ MODES_NUM (IN) : VECTOR (NM) INDICATING THE NUMBER OF HARMONICS OF EACH MODE
!!$ U_F_MODES (IN) : TRANSFORMATION, DIMENSION (D,D)
!!$ E_MULTIFLOQUET (IN) : MULTIMODE FLOQUET SPECTRUM
!!$ D_BARE (IN) : DIMENSION OF THE BARE HILBERT SPACE
!!$ FIELD (IN) : STRUCTURE DESCRIBING THE COUPLINGS
!!$ T1 (IN) : INITIAL TIME
!!$ T2 (IN) : FINAL TIME
!!$ U (OUT) : TRANSFORMATION BETWEEN THE EXTENDED BARE BASIS AND
!!$ THE FLOQUET STATES, DIMENSION (D_BARE,D)
!!$ INFO (INOUT): (POSSIBLE) ERROR FLAG

```

```

USE TYPES
USE SUBINTERFACE_LAPACK

```

```

IMPLICIT NONE

```

```

INTEGER, INTENT(IN) :: D,D_BARE,NM
INTEGER, INTENT(INOUT) :: INFO
INTEGER, DIMENSION(NM), INTENT(IN) :: MODES_NUM
TYPE(MODE), DIMENSION(NM), INTENT(IN) :: FIELD
DOUBLE PRECISION, INTENT(IN) :: T1,T2
DOUBLE PRECISION, DIMENSION(D), INTENT(IN) :: E_MULTIFLOQUET
COMPLEX*16, DIMENSION(D,D), INTENT(IN) :: U_F_MODES
COMPLEX*16, DIMENSION(D_BARE,D_BARE), INTENT(OUT) :: U

```

```

SUBROUTINE MULTIMODEFLOQUETMATRIX(ATOM_,NM,NF,MODES_NUM,FIELD,INFO)

```

```

!ID,size(modes_num,1),total_frequencies,MODES_NUM,FIELDS,INFO
! USE FLOQUET
!ATOM_ type atom, -> dimension of the bare Hilbert space
!NM -> number of modes
!NF -> Number of Fields
!MODES_NUM -> number of harmonics of each mode
!FIELD -> Field couplings
!INFO

```

```

USE ARRAYS
USE ATOMIC_PROPERTIES
USE TYPES
USE SUBINTERFACE_LAPACK

```

```

IMPLICIT NONE

```

```

INTEGER, INTENT(IN) :: NM,NF
INTEGER, INTENT(INOUT) :: INFO
INTEGER, DIMENSION(NM), INTENT(IN) :: MODES_NUM
TYPE(MODE), DIMENSION(NF), INTENT(IN) :: FIELD
TYPE(ATOM), INTENT(IN) :: ATOM_

```

```

SUBROUTINE MULTIMODEFLOQUETMATRIX_SP(ATOM_,NM,NF,MODES_NUM,FIELDS,VALUES_,ROW_INDEX_,COLUMN_,INFO)

```

```

!ATOM_      (IN)      : type of quantum system
!NM          (IN)      : number of modes
!NF          (IN)      : number of driving fields
!MODES_NUM   (IN)      : vector indicating the number of harmonics of each driving field (mode)
!FIELDS      (IN)      : Fields
!VALUES_     (OUT)     : Hamiltonian values
!ROW_INDEX_  (OUT)     : vector indicating the row position of values
!COLUMN_     (OUT)     : vector indicating the column position of the values
!INFO        (INOUT)   : error flag. INFO=0 means there is no error

```

```

USE TYPES          !(modes.f90)
USE MERGINGARRAYS !(utils.f90)

```

```

IMPLICIT NONE

```

```

INTEGER,          ,          INTENT(IN)      :: NM,NF
TYPE(MODE), DIMENSION(NF),    INTENT(INOUT)  :: FIELDS
TYPE(ATOM),       ,          INTENT(IN)      :: ATOM_
INTEGER, DIMENSION(NM),       INTENT(IN)      :: MODES_NUM
INTEGER,          ,          INTENT(INOUT)   :: INFO
COMPLEX*16, DIMENSION(:), ALLOCATABLE, INTENT(OUT) :: VALUES_
INTEGER, DIMENSION(:), ALLOCATABLE, INTENT(OUT) :: COLUMN_
INTEGER, DIMENSION(:), ALLOCATABLE, INTENT(OUT) :: ROW_INDEX_

```

```

SUBROUTINE MULTIMODEFLOQUETTRANSFORMATION(D,NM,MODES_NUM,U_F_MODES,E_MULTIFLOQUET,
                                           D_BARE,FIELD,T1,U,INFO)

```

```

! TIME-DEPENDENT TRANSFORMATION BETWEEN THE EXTENDED BARE BASIS AND THE FLOQUET STATES
!  $U(T1) = \sum_n U^n \exp(i n \omega T1)$ 
!

```

```

!!$ D          (IN)      : DIMENSION OF THE EXTENDED HILBERT SPACE (SIZE OF THE MULTIMODE FLOQUET)
!!$ NM         (IN)      : NUMBER OF MODES
!!$ MODES_NUM   (IN)      : VECTOR (NM) INDICATING THE NUMBER OF HARMONICS OF EACH MODE
!!$ U_F_MODES   (IN)      : TRANSFORMATION, DIMENSOON (D,D)
!!$ E_MULTIFLOQUET (IN)    : MULTIMODE FLOQUET SPECTRUM
!!$ D_BARE      (IN)      : DIMENSION OF THE BARE HILBERT SPACE
!!$ FIELD       (IN)      : STRUCTURE DESCRIBING THE COUPLINGS
!!$ T1          (IN)      : TIME. THE BARE 2 DRESSED TRANSFORMATINO IS TIME DEPENDENT
!!$ U           (OUT)     : TRANFORMATION BETWEEN THE EXTENDED BARE BASIS AND THE FLOQUET STATES,
!!$              DIMENSION (D_BARE,D)
!!$ INFO        (INOUT): (POSSIBLE) ERROR FLAG

```

```

USE TYPES

```

```

IMPLICIT NONE

```

```

INTEGER,          ,          INTENT(IN)      :: D,D_BARE,NM
INTEGER,          ,          INTENT(INOUT)   :: INFO
INTEGER, DIMENSION(NM),       INTENT(IN)      :: MODES_NUM
TYPE(MODE), DIMENSION(NM),    INTENT(IN)      :: FIELD
DOUBLE PRECISION,          ,   INTENT(IN)      :: T1
DOUBLE PRECISION, DIMENSION(D), INTENT(IN)      :: E_MULTIFLOQUET
COMPLEX*16, DIMENSION(D,D),    INTENT(IN)      :: U_F_MODES
COMPLEX*16, DIMENSION(D_BARE,D), INTENT(OUT)   :: U

```

```

SUBROUTINE MULTIMODEMICROMOTION(ID,D,NM,MODES_NUM,U_F_MODES,E_MULTIFLOQUET,D_BARE,FIELD,T1,U,INFO)

```

```

! TIME-DEPENDENT TRANSFORMATION BETWEEN THE EXTENDED BARE BASIS AND THE FLOQUET STATES
!  $U(T1) = \sum U^n \exp(i n \omega T1)$ 
!
!!$ D (IN) : DIMENSION OF THE EXTENDED HILBERT SPACE (SIZE OF THE MULTIMODE FLOQUET
!!$ NM (IN) : NUMBER OF MODES
!!$ MODES_NUM (IN) : VECTOR (NM) INDICATING THE NUMBER OF HARMONICS OF EACH MODE
!!$ U_F_MODES (IN) : TRANSFORMATION, DIMENSOON (D,D)
!!$ E_MULTIFLOQUET (IN) : MULTIMODE FLOQUET SPECTRUM
!!$ D_BARE (IN) : DIMENSION OF THE BARE HILBERT SPACE
!!$ FIELD (IN) : STRUCTURE DESCRIBING THE COUPLINGS
!!$ T1 (IN) : TIME. THE BARE 2 DRESSED TRANSFORMATINO IS TIME DEPENDENT
!!$ U (OUT) : TRANFORMATION BETWEEN THE EXTENDED BARE BASIS AND THE FLOQUET STATES, D
!!$ INFO (INOUT): (POSSIBLE) ERROR FLAG

```

```

!USE TYPES_C
USE TYPES
!USE MODES_4F
USE SUBINTERFACE_LAPACK
USE ATOMIC_PROPERTIES

```

```

IMPLICIT NONE

```

```

TYPE(ATOM),          INTENT(IN)      :: ID
INTEGER,              INTENT(IN)      :: D,D_BARE,NM
INTEGER,              INTENT(INOUT)   :: INFO
INTEGER,              DIMENSION(NM),  INTENT(IN)      :: MODES_NUM
TYPE(MODE),           DIMENSION(NM),  INTENT(IN)      :: FIELD
DOUBLE PRECISION,     INTENT(IN)      :: T1
DOUBLE PRECISION, DIMENSION(D),       INTENT(IN)      :: E_MULTIFLOQUET
COMPLEX*16,           DIMENSION(D,D),  INTENT(IN)      :: U_F_MODES
COMPLEX*16,           DIMENSION(D_BARE,D_BARE), INTENT(OUT) :: U

```

```

SUBROUTINE MICROMOTIONFOURIERDRESSED BASIS(ID,DRESSINGFIELDS_INDICES,MODES_NUM,FIELDS,
                                           U_FD,E_DRESSED,INFO)

```

```

! ID (in) :: TYPE(ATOM) system ID
! DRESSINGFIELDS_INDICES (in) :: integer array indicating the indices of the dressing modes
! MODES_NUM (in) :: integer array indicating the number of harmonics of all driving modes
! FIELDS (in) :: Array of TYPE(MODE) of dimension
! U_FD (out) :: complex*16 matrix fourier decomposition of the micromotion operator of the dr
! E_DRESSED (out) :: dressed energies
! INFO (inout) :: error flag
USE TYPES

```

```

TYPE(ATOM),          INTENT(IN) :: ID
INTEGER, DIMENSION(:), INTENT(IN) :: DRESSINGFIELDS_INDICES
INTEGER, DIMENSION(:), INTENT(IN) :: MODES_NUM
TYPE(MODE), DIMENSION(:), INTENT(IN) :: FIELDS
COMPLEX*16, DIMENSION(:,:), ALLOCATABLE, INTENT(OUT) :: U_FD
DOUBLE PRECISION, DIMENSION(:), ALLOCATABLE, INTENT(OUT) :: E_DRESSED

```

```

SUBROUTINE MICROMOTIONDRESSED BASIS(ID,MODES_NUM,DRESSINGFIELDS_INDICES,FIELDS,
                                     U_F_MODES,E_MULTIFLOQUET,T1,U,INFO)

```

```

! ID (in) :: TYPE(ATOM) system ID
! MODES_NUM (in) :: integer array indicating the number of harmonics of each driving mode
! DRESSINFIELDS_INDICES :: integer array indicating the indices of the dressing modes

```

```

! FIELDS          :: Array of TYPE(MODES) with NM components (all driving fields)
! U_F_MODES       :: complex*16 matrix of dimension Dx D. Fourier decomposition of the micromotion operator
! E_MULTIFLOQUET :: dressed energies
! T1              :: double precision, time
! U               :: complex*16 matrix of dimension D_BARE x D_BARE. micromotion operator at time T1
! INFO            :: error flag

```

```

USE TYPES

```

```

IMPLICIT NONE

```

```

TYPE(ATOM),          INTENT(IN)      :: ID
INTEGER,             DIMENSION(:),   INTENT(IN)  :: MODES_NUM
INTEGER,             DIMENSION(:),   INTENT(IN)  :: DRESSINGFIELDS_INDICES
COMPLEX*16,          DIMENSION(:,:), INTENT(IN)  :: U_F_MODES
DOUBLE PRECISION,    DIMENSION(:),   INTENT(IN)  :: E_MULTIFLOQUET
TYPE(MODE),          DIMENSION(:),   INTENT(IN)  :: FIELDS
DOUBLE PRECISION,    ,               INTENT(IN)  :: T1
COMPLEX*16,          DIMENSION(:,:), INTENT(OUT) :: U
INTEGER,             INTENT(INOUT)   :: INFO

```

```

SUBROUTINE MULTIMODETRANSITIONAVG(D,NM,FIELD,MODES_NUM,U_F_MODES,
                                   E_MULTIFLOQUET,D_BARE,U,INFO)

```

```

!!$  AVERAGE TIME EVOLUTION OPERATOR OF A MULTIMODE DRESSED SYSTEM.
!!$  THE AVERAGE EVOLUTION OPERATOR IS WRITTEN IN THE BASIS USED TO EXPRESS THE
!!$  MULTIMODE FLOQUET HAMILTONIAN
!!$  U : MATRIX OF AVERAGE TRANSITION PROBABILITIES
!!$
!!$  D          (IN)   : DIMENSION OF THE EXTENDED HILBERT SPACE
!!$                (SIZE OF THE MULTIMODE FLOQUET MATRIX)
!!$  NM         (IN)   : NUMBER OF MODES
!!$  MODES_NUM  (IN)   : VECTOR (NM) INDICATING THE NUMBER OF HARMONICS OF EACH MODE
!!$  U_F_MODES  (IN)   : TRANSFORMATION, DIMENSION (D,D)
!!$  E_MULTIFLOQUET (IN) : MULTIMODE FLOQUET SPECTRUM
!!$  D_BARE     (IN)   : DIMENSION OF THE BARE HILBERT SPACE
!!$  U          (OUT)  : MATRIX OF AVERAGE TRANSITION PROBABILITIES
!!$  INFO      (INOUT): (POSSIBLE) ERROR FLAG

```

```

USE TYPES

```

```

IMPLICIT NONE

```

```

TYPE(MODE),DIMENSION(NM), INTENT(IN)      :: FIELD
INTEGER,    DIMENSION(NM), INTENT(IN)     :: MODES_NUM

INTEGER,          INTENT(IN)      :: D,D_BARE,NM
INTEGER,          INTENT(INOUT)   :: INFO
DOUBLE PRECISION, DIMENSION(D),   INTENT(IN)  :: E_MULTIFLOQUET
COMPLEX*16,       DIMENSION(D,D),  INTENT(IN)  :: U_F_MODES
DOUBLE PRECISION, DIMENSION(D_BARE,D_BARE), INTENT(OUT) :: U

```

14 DRIVER SUBROUTINES

```
SUBROUTINE DRESSED BASIS(D, ID, NM, MODES_NUM, FIELDS, U_FD, E_DRESSED, INFO)
```

```
!!$ THIS SUBROUTINES CALCULATES THE FOURIER COMPONENTS OF THE
!!$ TRANSFORMATION BETWEEN THE BARE BASIS TO THE DRESSED BASIS DEFINED
!!$ BY THE FULL SET OF DRIVING FIELDS.
!!$
!!$ D : DIMENSION OF THE MULTIMODE EXTENDED HILBERT SPACE
!!$ ID (IN) : TYPE OF QUANTUM SYSTEM
!!$ NM (IN) : NUMBER OF MODES == NUMBER OF DRIVING FIELDS
!!$ MODES_NUM : VECTOR INDICATING THE NUMBER OF HARMONICS OF EACH DRESSING FIELD
!!$ FIELDS (IN) : AMPLITUDE, FREQUENCY AND PHASES OF ALL DRIVING FIELDS
!!$ U_FD (OUT) : THIS IS THE TRANSFORMATION WE ARE LOOKING FOR
!!$ E_DRESSED (OUT) : DRESSED ENERGIES
!!$ INFO (INOUT) : INFO = 0 MEANS SUCESS
```

```
USE ATOMIC_PROPERTIES
USE TYPES
USE SUBINTERFACE
USE SUBINTERFACE_LAPACK
USE FLOQUETINIT_
USE ARRAYS
```

```
IMPLICIT NONE
TYPE(MODE), DIMENSION(NM), INTENT(IN) :: FIELDS
TYPE(ATOM), INTENT(IN) :: ID
INTEGER, DIMENSION(NM), INTENT(IN) :: MODES_NUM
COMPLEX*16, DIMENSION(D,D), INTENT(OUT) :: U_FD
DOUBLE PRECISION, DIMENSION(D), INTENT(OUT) :: E_DRESSED
INTEGER, INTENT(IN) :: NM, D
INTEGER, INTENT(INOUT) :: INFO
```

```
SUBROUTINE DRESSED BASIS_SP(D, ID, NM, MODES_NUM, FIELDS, U_FD, E_DRESSED, INFO)
```

```
!!$ THIS SUBROUTINES CALCULATES THE TRANSFORMATION BETWEEN THE BARE
!!$ BASIS TO THE DRESSED BASIS DEFINED BY THE FULL SET OF DRIVING FIELDS.
!!$ D : DIMENSION OF THE MULTIMODE EXTENDED HILBERT SPACE
!!$ ID (IN) : TYPE OF QUANTUM SYSTEM
!!$ NM (IN) : NUMBER OF MODES == NUMBER OF DRIVING FIELDS
!!$ MODES_NUM : VECTOR INDICATING THE NUMBER OF HARMONICS OF EACH DRESSING FIELD
!!$ FIELDS (IN) : AMPLITUDE, FREQUENCY AND PHASES OF ALL DRIVING FIELDS
!!$ U_FD (OUT) : THIS IS THE TRANSFORMATION WE ARE LOOKING FOR
!!$ E_DRESSED (OUT) : DRESSED ENERGIES
!!$ INFO (INOUT) : INFO = 0 MEANS SUCESS
```

```
USE ATOMIC_PROPERTIES
USE TYPES
USE SPARSE_INTERFACE
USE SUBINTERFACE
USE SUBINTERFACE_LAPACK
USE FLOQUETINIT_
USE ARRAYS
```

```
IMPLICIT NONE
TYPE(MODE), DIMENSION(NM), INTENT(INOUT) :: FIELDS
```



```

TYPE(ATOM),                                INTENT(IN)      :: ID
INTEGER,    DIMENSION(NM),                 INTENT(IN)      :: MODES_NUM
COMPLEX*16, DIMENSION(D,D),                 INTENT(OUT)     :: U_FD
DOUBLE PRECISION, DIMENSION(D), INTENT(OUT)  :: E_DRESSED
INTEGER,                                INTENT(IN)      :: NM,D
INTEGER,                                INTENT(INOUT)   :: INFO

```

```

SUBROUTINE TIMEEVOLUTIONOPERATOR(ID,D_BARE,NM,MODES_NUM,FIELD,T1,T2,U,INFO)
! TIME EVOLUTION OPERATOR OF A MULTIMODE DRESSED SYSTEM. THE EVOLUTION
! OPERATOR IS WRITTEN IN THE BASIS USED TO EXPRESS THE
! MULTIMODE FLOQUET HAMILTONIAN
! U : MATRIX OF AMPLITUDES OF PROBABILITIES FOR TRANSITIONS BETWEEN T1 TO T2
!!$ NM          (IN)      : NUMBER OF MODES
!!$ MODES_NUM    (IN)      : VECTOR (NM) INDICATING THE NUMBER OF HARMONICS OF EACH MODE
!!$ D_BARE       (IN)      : DIMENSION OF THE BARE HILBERT SPACE
!!$ FIELD        (IN)      : STRUCTURE DESCRIBING THE COUPLINGS
!!$ T1           (IN)      : INITIAL TIME
!!$ T2           (IN)      : FINAL TIME
!!$ U            (OUT)     : TRANSFORMATION BETWEEN THE EXTENDED BARE BASIS AND
!!$              THE FLOQUET STATES, DIMENSION (D_BARE,D)
!!$ INFO         (INOUT): (POSSIBLE) ERROR FLAG

```

```

USE ATOMIC_PROPERTIES
USE TYPES
USE SUBINTERFACE
USE SUBINTERFACE_LAPACK
USE FLOQUETINIT_
USE ARRAYS

```

```

IMPLICIT NONE
TYPE(ATOM) ,                                INTENT(IN)      :: ID
INTEGER,                                INTENT(IN)      :: D_BARE
INTEGER,                                INTENT(IN)      :: NM
INTEGER,    DIMENSION(NM),                 INTENT(IN)      :: MODES_NUM
TYPE(MODE),    DIMENSION(NM),                 INTENT(IN)      :: FIELD
DOUBLE PRECISION,                                INTENT(IN)      :: T1
DOUBLE PRECISION,                                INTENT(IN)      :: T2
COMPLEX*16,    DIMENSION(D_BARE,D_BARE), INTENT(OUT)     :: U
INTEGER,                                INTENT(INOUT)   :: INFO

```

```

SUBROUTINE MICROMOTIONFOURIERDRESSED BASIS(ID,DRESSINGFIELDS_INDICES,
                                           MODES_NUM,FIELDS, U_FD,E_DRESSED,INFO)
! THIS SUBROUTINE CALCULATES THE FOURIER COMPONENTS (U_FD) AND PHASES (E_DRESSED)
! OF THE MICROMOTION OPERATOR OF SUBSET OF DRIVING MODES
! ID      (in)      :: TYPE(ATOM) system ID
! DRESSINGFIELDS_INDICES (in) :: integer array indicating the indices of the dressing modes
! MODES_NUM (in)      :: integer array indicating the number of harmonics of all driving modes
! FIELDS    (in)      :: Array of TYPE(MODE) of dimension
! U_FD      (out)     :: complex*16 matrix fourier decomposition of the micromotion
!              operator of the dressed basis
! E_DRESSED (out)     :: dressed energies
! INFO      (inout)  :: error flag

```

```

USE TYPES
IMPLICIT NONE
TYPE(ATOM),                                INTENT(IN)      :: ID
INTEGER,    DIMENSION(:),                 INTENT(IN)      :: DRESSINGFIELDS_INDICES
INTEGER,    DIMENSION(:),                 INTENT(IN)      :: MODES_NUM

```

```

TYPE(MODE), DIMENSION(:),          INTENT(IN)  :: FIELDS
COMPLEX*16, DIMENSION(:,:),        ALLOCATABLE, INTENT(OUT) :: U_FD
DOUBLE PRECISION, DIMENSION(:),    ALLOCATABLE, INTENT(OUT) :: E_DRESSED
INTEGER, INTENT(INOUT) :: INFO

```

```

END SUBROUTINE MICROMOTIONFOURIERDRESSED BASIS

```

```

SUBROUTINE MICROMOTIONDRESSED BASIS(ID, MODES_NUM, DRESSINGFIELDS_INDICES, FIELDS,
                                     U_F_MODES, E_MULTIFLOQUET, T1, U, INFO)

```

```

! THIS SUBROUTINE CALCULATES U: THE TIME-DEPENDENT MICROMOTION OPERATOR OF
! A SUBSET OF THE DRIVING MODES. U_F_MODES AND E_MULTIFLOQUET ARE THE ARRAYS
! CALCULATED WITH THE SUBROUTINE MICROMOTIONFOURIERDRESSED BASIS

```

```

! ID (in)          :: TYPE(ATOM) system ID
! MODES_NUM (in)  :: integer array indicating the number of harmonics of each driving mode
! DRESSINGFIELDS_INDICES :: integer array indicating the indices of the dressing modes
! FIELDS          :: Array of TYPE(MODES) with NM components (all driving fields)
! U_F_MODES       :: complex*16 matrix of dimension DxD. Fourier decomposition of
!                  the micromotion operator of the dressed basis
! E_MULTIFLOQUET :: dressed energies
! T1              :: double precision, time
! U               :: complex*16 matrix of dimension D_BARE x D_BARE. micromotion operator
!                  at time T1
! INFO            :: error flag

```

```

USE TYPES
IMPLICIT NONE

```

```

TYPE(ATOM),          INTENT(IN)  :: ID
INTEGER,             DIMENSION(:), INTENT(IN)  :: MODES_NUM
INTEGER,             DIMENSION(:), INTENT(IN)  :: DRESSINGFIELDS_INDICES
COMPLEX*16,          DIMENSION(:,:), INTENT(IN)  :: U_F_MODES
DOUBLE PRECISION,    DIMENSION(:),  INTENT(IN)  :: E_MULTIFLOQUET
TYPE(MODE),          DIMENSION(:),  INTENT(IN)  :: FIELDS
DOUBLE PRECISION ,   INTENT(IN)  :: T1
COMPLEX*16,          DIMENSION(:,:), INTENT(OUT) :: U
INTEGER,             INTENT(INOUT) :: INFO

```

14.1 Utility subroutines

```
SUBROUTINE PACKINGBANDMATRIX(N,A,KD,AB,INFO)
```

```
! brute force packing of a banded matrix
```

```
IMPLICIT NONE
INTEGER, INTENT(INOUT) :: INFO
INTEGER, INTENT(IN)    :: N,KD
COMPLEX*16, DIMENSION(N,N) :: A
COMPLEX*16, DIMENSION(KD+1,N) :: AB
```

```
SUBROUTINE LAPACK_FULLEIGENVALUES(H,N,W_SPACE,INFO)
```

```
!eigenvalues/vectors of matrix ab
```

```
!H, inout, packed banded matrix
```

```
! , out,eigenvectors
```

```
!N, in,matrix dimension
```

```
!W_space, out, eigenvalues
```

```
!INFO,inout, error flag
```

```
!H is COMPLEX*16 array, dimension (N, N)
```

```
! 69 *>      On entry, the Hermitian matrix A.  If UPLO = 'U', the
```

```
! 70 *>      leading N-by-N upper triangular part of A contains the
```

```
! 71 *>      upper triangular part of the matrix A.  If UPLO = 'L',
```

```
! 72 *>      the leading N-by-N lower triangular part of A contains
```

```
! 73 *>      the lower triangular part of the matrix A.
```

```
! 74 *>      On exit, if JOBZ = 'V', then if INFO = 0, A contains the
```

```
! 75 *>      orthonormal eigenvectors of the matrix A.
```

```
! 76 *>      If JOBZ = 'N', then on exit the lower triangle (if UPLO='L')
```

```
! 77 *>      or the upper triangle (if UPLO='U') of A, including the
```

```
! 78 *>      diagonal, is destroyed.
```

```
!
```

```
! The eigenvector H(:,r) corresponds to the eigenvalue W_SPACE(r)
```

```
!
```

```
IMPLICIT NONE
```

```
INTEGER,                                INTENT(IN)    :: N
```

```
COMPLEX*16,          DIMENSION(N,N), INTENT(INOUT) :: H
```

```
DOUBLE PRECISION, DIMENSION(N),  INTENT(INOUT) :: W_SPACE
```

```
INTEGER,                                INTENT(OUT)   :: INFO
```

```
SUBROUTINE LAPACK_FULLEIGENVALUESBAND(AB,Z,KD,N,W,INFO)
```

```
!eigenvalues/vectors of banded matrix ab
```

```
!AB, inout, packed banded matrix
```

```
!Z, out,eigenvectors
```

```
!KD out, calcuated eigenvectors
```

```
!N, in,matrix dimension
```

```
!W, out, eigenvalues
```

```
!INFO,inout, error flag
```

```
!H is COMPLEX*16 array, dimension (N, N)
```

```
! 69 *>      On entry, the Hermitian matrix A.  If UPLO = 'U', the
```

```
! 70 *>      leading N-by-N upper triangular part of A contains the
```

```
! 71 *>      upper triangular part of the matrix A.  If UPLO = 'L',
```

```
! 72 *>      the leading N-by-N lower triangular part of A contains
```

```
! 73 *>      the lower triangular part of the matrix A.
```

```
! 74 *>      On exit, if JOBZ = 'V', then if INFO = 0, A contains the
```

```
! 75 *>      orthonormal eigenvectors of the matrix A.
```

```
! 76 *>      If JOBZ = 'N', then on exit the lower triangle (if UPLO='L')
```

```
! 77 *>      or the upper triangle (if UPLO='U') of A, including the
```

```

! 78 *>          diagonal, is destroyed.
!
! The eigenvector H(:,r) corresponds to the eigenvalue W_SPACE(r)
!
IMPLICIT NONE
INTEGER,                                INTENT(IN)      :: N,KD
COMPLEX*16,          DIMENSION(KD+1,N), INTENT(INOUT)   :: AB
COMPLEX*16,          DIMENSION(N,N),    INTENT(INOUT)   :: Z
DOUBLE PRECISION,    DIMENSION(N),      INTENT(INOUT)   :: W
INTEGER,                                INTENT(OUT)     :: INFO

```

```

SUBROUTINE LAPACK_SELECTEIGENVALUES(H,N,W_SPACE,L1,L2,Z,INFO)

```

```

!selected eigenvalues/vectors of hermitian matrix

```

```

!H, inout, packed banded matrix

```

```

! , out,eigenvectors

```

```

!N, in,matrix dimension

```

```

!W_space, out, eigenvalues

```

```

!L1 ordinal lowest eigenvalue

```

```

!L2 ordinal highest eigenvlaue

```

```

!Z : eigenvectors

```

```

!INFO,inout, error flag

```

```

!USE FLOQUET

```

```

IMPLICIT NONE

```

```

INTEGER,                                INTENT(IN)      :: N,L1,L2

```

```

COMPLEX*16, DIMENSION(:,:),            INTENT(INOUT)   :: H

```

```

COMPLEX*16, DIMENSION(:,:),            INTENT(OUT)     :: Z

```

```

DOUBLE PRECISION, DIMENSION(:),         INTENT(OUT)     :: W_SPACE

```

```

INTEGER,                                INTENT(OUT)     :: INFO

```

```

SUBROUTINE MKLSPARSE_FULLEIGENVALUES(D,DV,VALUES,ROW_INDEX,COLUMN,E_L,E_R,E_FLOQUET,U_F,INFO)

```

```

!CALCULATES THE ENERGY SPECTRUM OF THE MATRIX REPRESENTED BY VALUES, ROW_INDEX AND COLUMN

```

```

! D (IN), MATRIX DIMENSION == NUMBER OF EIGENVALUES

```

```

! DV (IN), NUMBER OF VALUES != 0

```

```

! VALUES (IN) ARRAY OF VALUES

```

```

! ROW_INDEX (IN), ARRAY OF INDICES

```

```

! COLUMN (IN),      ARRAY OF COLUMN NUMBERS

```

```

! E_L (IN),        LEFT BOUNDARY OF THE SEARCH INTERVAL

```

```

! E_R (IN),        RIGHT BOUNDARY OF THE SEARCH INTERVAL

```

```

! E_FLOQUET (OUT), ARRAY OF EIGENVALUES

```

```

! INFO      (INOUT) ERROR FLAG and VERBOSITY FLAG

```

```

!           0 display no information

```

```

!           1 DISPLAY INFORMAITON ABOUT THE SIZE OF THE ARRAYS

```

```

!           10 DISPLAY INFORMAITON ABOUT THE ARRAYS AND THE ARRAYS

```

```

USE FEAST

```

```

IMPLICIT NONE

```

```

INTEGER,                                INTENT(IN)      :: D,DV

```

```

COMPLEX*16,          DIMENSION(DV),    INTENT(INOUT)   :: VALUES

```

```

INTEGER,            DIMENSION(DV),     INTENT(INOUT)   :: COLUMN

```

```

INTEGER,            DIMENSION(D+1),    INTENT(INOUT)   :: ROW_INDEX

```

```

DOUBLE PRECISION,          INTENT(IN)   :: E_L,E_R

```

```

DOUBLE PRECISION, DIMENSION(D),        INTENT(OUT)     :: E_FLOQUET

```

```

COMPLEX*16,            DIMENSION(D,D), INTENT(OUT)     :: U_F

```

```

INTEGER,                                INTENT(INOUT)   :: INFO

```

```

SUBROUTINE QUICK_SORT_INTEGERS(v,index_t,N)

  IMPLICIT NONE
  INTEGER, INTENT(IN) :: N
  INTEGER, DIMENSION(N),INTENT(INOUT) :: v
  INTEGER, DIMENSION(N),INTENT(INOUT) :: index_t

  INTEGER, PARAMETER :: NN=10000, NSTACK=8000

```

```

SUBROUTINE WRITE_MATRIX(A)
! it writes a matrix of doubles nxm on the screen
  DOUBLE PRECISION, DIMENSION(:,:) :: A
  CHARACTER(LEN=105) STRING
  CHARACTER(LEN=105) aux_char
  integer :: aux

```

```

SUBROUTINE WRITE_MATRIX_INT(A)
!it writes a matrix of integer nxm on the screen
  INTEGER, DIMENSION(:,:) :: A

```

```

SUBROUTINE COORDINATEPACKING(D,A,V,R,C,index,INFO)
  IMPLICIT NONE
  INTEGER,INTENT(IN):: D
  COMPLEX*16,DIMENSION(D,D),INTENT(IN)  :: A
  COMPLEX*16,DIMENSION(D*D),INTENT(OUT) :: V
  INTEGER, DIMENSION(D*D), INTENT(OUT) :: R,C
  INTEGER, INTENT(OUT)  :: index
  INTEGER, INTENT(INOUT) :: INFO

```

```

SUBROUTINE APPENDARRAYS(V,B,INFO)
  COMPLEX*16, DIMENSION(:),ALLOCATABLE, INTENT(INOUT) :: V
  COMPLEX*16, DIMENSION(:),INTENT(IN)  :: B
  INTEGER,          INTENT(INOUT) :: INFO

```

```

SUBROUTINE APPENDARRAYSI(V,B,INFO)
  INTEGER, DIMENSION(:),ALLOCATABLE, INTENT(INOUT) :: V
  INTEGER, DIMENSION(:),INTENT(IN)  :: B
  INTEGER,          INTENT(INOUT) :: INFO

```

SUBROUTINE VARCRCPACKING(N,DIM,UPL0,zero,A,VALUES,COLUMNS,ROWINDEX,INFO)

INTEGER, INTENT(IN) :: N
INTEGER, INTENT(INOUT) :: INFO,DIM
CHARACTER, INTENT(IN) :: UPL0
DOUBLE PRECISION, INTENT(IN) :: ZERO
COMPLEX*16,DIMENSION(N,N), INTENT(IN) :: A

COMPLEX*16, DIMENSION(DIM), INTENT(OUT) :: VALUES
INTEGER, DIMENSION(DIM), INTENT(OUT) :: COLUMNS
INTEGER, DIMENSION(N+1), INTENT(OUT) :: ROWINDEX

15 C++ Wrappers prototypes: src/MultimodeFloquet.h

```
struct mode_c{
    double omega;
    dcmplx x,y,z;
    double phi_x,phi_y,phi_z;
    int N_Floquet;
};

struct atom_c{
    int id_system;
    int d_bare;
};

extern "C" {

    // DIMENSION OF THE MULTIMODE FLOQUET MATRIX. CALCULATED INTERNALLY
    int h_floquet_size;
    int h_floquet_c; // Floquet matrix in the bare basis

    // GENERAL INIT SUBROUTINE
    void floquetinit_qubit_c_ (atom_c *id, int *lenght_name, char * atomicspecie,
    void floquetinit_spin_c_ (atom_c *id, int *lenght_name, char * atomicspecie, double * jtotal,
    void floquetinit_alkali_c_(atom_c *id, int *lenght_name, char * atomicspecie, int * lenght_name2,

    // SET HAMILTONIAN OF SPIN-LIKE MODELS
    void sethamiltoniancomponents_c_(atom_c *id,int * nm, int * total_frequencies,int * modes_num,mod

    // BUILDING FLOQUET MATRIX OF GENERIC MODEL
    void multimodefloquetmatrix_c_ (atom_c *id,int * nm, int * total_frequencies,
        int * modes_num,mode_c * fields, int * info);
    void get_h_floquet_c_(int * h, dcmplx * values, int* info);
    int multimodefloquetmatrix_c_python_(atom_c *id,int * nm, int * total_frequencies,
        int * modes_num,mode_c * fields,int * info);
    void multimodefloquetmatrix_sp_c_ (atom_c *id,int * nm, int * total_frequencies,
        int * modes_num,mode_c * fields, int * info);
    void multimodefloquetmatrix_python_sp_c_(atom_c *id,int * nm, int * total_frequencies,
        int * modes_num,mode_c * fields, int * h_f,int * info);
    void get_h_floquet_sp_c_(int * h_f, dcmplx * values, int * row_index, int * column, int * info);

    // CALCULATE THE SPECTRUM OF THE FLOQUET HAMILTONIAN
    void lapack_fulleigenvalues_c_(dcmplx * u_f,int * h_floquet_size,
        double * e_floquet,int *info);
    void mklsparse_fulleigenvalues_c_(int * h_floquet_size,
        double * e_l,double * e_r,double * e_floquet,dcmplx *U_F, int * info);
    void matmul_c_(int * op, dcmplx * a, int * ra, int * ca,
        dcmplx * b, int * rb, int * cb, dcmplx * c,int * info);

    // CONTRSUTION OF THE TIME-EVOLUTION OPERATOR
    void multimodetransitionavg_c_(int * h_floquet_size,int * nm,mode_c * fields,
        int * modes_num,dcmplx * U_F,double * e_floquet,int * d_bare,double * p_avg,int *info);
    void multimodefloquettransformation_c_(int * h_floquet_size,int * nm,int * modes_num,
        dcmplx * U_F,double * e_floquet,int * d_bare,mode_c * fields,double * t1,dcmplx * U_B2D,int *
    void multimodemicrotation_c_(atom_c *id,int * h_floquet_size,int * nm,int * modes_num,
        dcmplx * U_F,double * e_floquet,int * d_bare,mode_c * fields,double * t1,dcmplx * U_B2D,int *
    void multimodetimeevolutionoperator_c_(int * h_floquet_size,int * nm,int * modes_num,
        dcmplx * U_F,double * e_floquet,int * d_bare,mode_c * fields,double * t1,double * t2,dcmplx *
```

```

void timeevolutionoperator_c(atom_c *id, int *d_bare, int *nm, int *nf, int * modes_num,
    mode_c *field, double *t1, double *t2, dcmplx *U, int *info);

// DEFINITION OF DRESSED BASIS
void      dressedbasis_c(int * h_floquet_size,atom_c *id,int * nm,
                        int * modes_num,mode_c * fields, dcmplx * U_FD,
                        double * e_dressed,int * info);
void  dressedbasis_subset_c(atom_c *id , int * dressingfloquetdimension,
                        int * dressingfields, int * nm, int * dressingfields_indices,
                        int * modes_num,mode_c * fields, dcmplx * U_FD,
                        double * e_dressed,int * info);
void  dressedbasis_subset_sp_c(atom_c * id, int * dressingfloquetdimension,
                        int * dressingfields,int * nm, int * dressingfields_indices,
                        int * modes_num,mode_c * fields, dcmplx * U_FD,
                        double * e_dressed,int * info);
void  dressedbasis_sp_c(int h_floquet_size, atom_c *id, int * nm, int * modes_num,
                        mode_c * fields, dcmplx * U_FD, double * e_dressed,
                        int * info);
void micromotionfourierdressedbasis_c(atom_c *id , int *DF, int * dressingfields_indices,
                        int * nm, int * modes_num, int *nf, mode_c * fields,
                        int * nd,dcmplx * U_FD, double *e_dressed,int * info);
void micromotiondressedbasis_c(atom_c *id , int * modes_num, int * dressingfields_indices,
                        mode_c * fields, double T1, dcmplx * U, int * info);

// UTILITY FUNCTIONS: WRITE MATRICES ON THE SCREEN
void write_matrix_c(double *A,int * A_dim);
void rec_write_matrix_c(double *A,int * A_dim1, int * A_dim2);

// deallocate all arrays allocated with fortran
void deallocateall_c(int *id);

```


16 Python Wrappers prototypes

OPENMMMF includes a set of wrappers to use the library with Python, defined in the file `src/openmmf.py`, The wrapper use CTYPES to pass on parameters to the C++ wrappers. In order for this to work, you should make sure that the openmmf dynamical library is loaded correctly, e.g. using the instruction: `openmmfC=ctypes.CDLL('../lib/libmultimodefloquet.so')`

near the top of `src/openmmf.py`. Make sure the path is correct.

The wrappers were developed and tested using: Python 3.7.6, ctypes 1.1.0, numpy 1.16.2, scipy 1.4.1 and matplotlib 3.0.3.

```
#####
#   // GENERAL CLASSES
#####

class atom_c_T(ctypes.Structure):
    _fields_ = [
        ("id_system", c_int),
        ("d_bare", c_int)
    ]

#####
#####

class mode_c_T(ctypes.Structure):
    c_dcplx = ctypes.c_double*2
    _fields_ = [
        ("omega", c_double),
        ("x", c_dcplx),
        ("y", c_dcplx),
        ("z", c_dcplx),
        ("phi_x", c_double),
        ("phi_y", c_double),
        ("phi_z", c_double),
        ("N_Floquet", c_int)
    ]

#####
#   // GENERAL INIT SUBROUTINE
#####

openmmf.floquetinit(id,*argsv,info):
    Parameters: id: atom_c_T
                Identifies the type of system
                *argsv: array
                Depending on the system of inters
                info: int
                Error/success flag

This function supports the following calls:

openmmf.floquetinit(id,qubit,info=info)
    qubit: str
            'qubit' to define a two-level system
    info: int
            Error/success flag

openmmf.floquetinit(id,alkali,manifold,info=info)
    alkali: str
```

```

        '87Rb', '23Na' atomic specie
manifold: str
    Manifold of the hyperfine splitting
    'L' : Lower with total angular moment  $F = I - J$ 
    'U' : Upper with total angular momentum  $F = I + J$ 
    'B' : both manifolds
info: int
    Error/success flag

openmmf.floquetinit(id,system,L,info=info)
    system: str
        'spin'    single particle with total angular momentum  $2*L$ 
        'lattice' quantum system with L sites/states
    L: np.double
        define
info: int
    Error/success flag

#=====
#  EVALUATE THE HAMILTONIAN COMPONENTS OF PRE-DEFINED SYSTEMS.
#=====
openmmf.sethamiltoniancomponents(id,modes_num,fields,info):
    id: atom_c_T instance
    modes_num: 1D np.array([],dtype=np.int32)
        define the number of fundamental frequencies and modes
        the total number of modes is nm = np.sum(modes_num)
    fields: instance of openmmf.mode_c_T*nm
        equivalent to the FORTRAN data derived type FIELDS
    info: int
        Error/success flag

#=====
#  // BUILDING FLOQUET MATRIX OF GENERIC MODEL
#=====
openmmf.multimodefloquetmatrix(id,modes_num,fields,info):
    id: atom_c_T instance
    modes_num: 1D np.array([],dtype=np.int32)
        define the number of fundamental frequencies and modes
        the total number of modes is nm = np.sum(modes_num)
    fields: instance of openmmf.mode_c_T*nm
        equivalent to the FORTRAN data derived type FIELDS
    info: int
        Error/success flag

openmmf.get_h_floquet(h_floquet_size, info):
    h_floquet_size: int
        size of the multimode Floquet matrix
    returns the multimode floquet matrix

openmmf.multimodefloquetmatrix_sp(id,modes_num,fields,info):
    id: atom_c_T instance
    modes_num: 1D np.array([],dtype=np.int32)
        define the number of fundamental frequencies and modes
        the total number of modes is nm = np.sum(modes_num)
    fields: instance of openmmf.mode_c_T*nm
        equivalent to the FORTRAN data derived type FIELDS
    info: int
        Error/success flag

```

```

openmmf.get_h_floquet_sp(h_floquet_size, info):
    h_floquet_size: int
        size of the multimode Floquet matrix
    returns VALUE,ROW,COLUMN arrays to represent a sparse multimode floquet matrix

#####
# // CALCULATE THE SPECTRUM OF THE FLOQUET HAMILTONIAN
#####
openmmf.lapack_fulleigenvalues(U_F,h_floquet_size,e_floquet,info):
    U_F: 1D np.array([],dtype=np.complex)
        h_floquet_size x h_floquet_size matrix of eigenvectors
    h_floquet_size: int
        size of the multimode Floquet matrix
    e_floquet: 1D np.array([],dtype=np.complex)
        array of size h_floquet_size containing the eigenvalues of
        the multimode Floquet Hamiltonian
    info: int
        Error/success flag

#####
# EVALUATES THE TIME AVERAGE TRANSITIONS PROBABILITIES
#####
openmmf.multimodetransitionavg(h_floquet_size,fields,modes_num,U_F,e_floquet,d_bare,p_avg,info):
    h_floquet_size: int
        size of the multimode Floquet matrix
    fields: instance of openmmf.mode_c_T*nm
        equivalent to the FORTRAN data derived type FIELDS
    modes_num: 1D np.array([],dtype=np.int32)
        define the number of fundamental frequencies and modes
        the total number of modes is nm = np.sum(modes_num)
    U_F: 1D np.array([],dtype=np.complex)
        h_floquet_size x h_floquet_size matrix of eigenvectors
    e_floquet: 1D np.array([],dtype=np.complex)
        array containing the eigenvalues of the multimode Floquet
        Hamiltonian
    d_bare: int
        Hilbert space dimension of the static system
    p_avg: 1D np.array([],dtype=np.double)
        d_bare x d_bare Matrix of the time-average state occupations
    info: int
        Error/success flag

#####
# FOURIER COMPONENTS OF THE TRANSFORMATION BETWEEN THE DRESSED AND THE BARE BASIS
# U_B2D is of dimension [D_bare*D_floquet]
#####
openmmf.multimodefloquettransformation(modes_num, U_F,e_floquet,d_bare,fields,t1, U_B2D,info):
    modes_num: 1D np.array([],dtype=np.int32)
        define the number of fundamental frequencies and modes
        the total number of modes is nm = np.sum(modes_num)
    U_F: 1D np.array([],dtype=np.complex)
        h_floquet_size x h_floquet_size Matrix of eigenvectors
    e_floquet: 1D np.array([],dtype=np.complex)
        array containing the eigenvalues of the multimode Floquet
        Hamiltonian
    d_bare: int
        Hilbert space dimension of the static system
    fields: instance of openmmf.mode_c_T*nm
        equivalent to the FORTRAN data derived type FIELDS
    t1: np.double
        instant of time to evaluate the micromotion operator
    U_B2D: 1D np.array([],dtype=np.complex)

```

```

        d_bare x d_bare Matrix of the time-average state occupationrs
info: int
    Error/success flag

=====
# CALCULATE THE MICROMOTION OPERATOR
# U_B2D is of dimension [D_bare*D_bare]
=====
openmmf.multimodemicromotion(id,h_floquet_size,modes_num,U_F,e_floquet,d_bare,fields,t1, U_B2D,info)
    id: atom_c_T instance
    h_floquet_size: int
        size of the multimode Floquet matrix
    modes_num: 1D np.array([],dtype=np.int32)
        define the number of fundamental frequencies and modes
        the total number of modes is nm = np.sum(modes_num)
    U_F: 1D np.array([],dtype=np.complex)
        h_floquet_size x h_floquet_size Matrix of eigenvectors
    e_floquet: 1D np.array([],dtype=np.complex)
        array containing the eigenvalues of the multimode Floquet
        Hamiltonian
    d_bare: int
        Hilbert space dimension of the static system
    fields: instance of openmmf.mode_c_T*nm
        equivalent to the FORTRAN data derived type FIELDS
    t1: np.double
        instant of time to evaluate the micromotion operator
    U_B2D: 1D np.array([],dtype=np.complex)
        d_bare x d_bare Matrix of the time-average state occupationrs
    info: int
        Error/success flag

=====
# COMP. ROUTINE: EVALUATE THE TIME EVOULUTION OPERATOR BETWEEN T1 AND T2
# USING THE FOURIER DECOMPOSITION
# OF THE MICROMOTION OPERATOR, WHICH IS STORE IN U_F.
=====
def multimodetimeevolutionoperator(h_floquet_size,modes_num,U_F,e_floquet,d_bare,
    fields,t1,t2,U_AUX,info):
    h_floquet_size: int
        size of the multimode Floquet matrix
    modes_num: 1D np.array([],dtype=np.int32)
        define the number of fundamental frequencies and modes
        the total number of modes is nm = np.sum(modes_num)
    U_F: 1D np.array([],dtype=np.complex)
        h_floquet_size x h_floquet_size Matrix of eigenvectors
    e_floquet: 1D np.array([],dtype=np.complex)
        array containing the eigenvalues of the multimode Floquet
        Hamiltonian
    d_bare: int
        Hilbert space dimension of the static system
    fields: instance of openmmf.mode_c_T*nm
        equivalent to the FORTRAN data derived type FIELDS
    t1: np.double
        instant of time to evaluate the time-evolution operator
    t2: np.double
        instant of time to evaluate the time-evolution operator
    U_AUX: 1D np.array([],dtype=np.complex)
        d_bare x d_bare Matrix of the time-evolution operator
    info: int
        Error/success flag

```

```

=====
# DRIVER ROUTINE: EVALUATE THE TIME EVOLUTION OPERATOR BETWEEN T1 AND T2
=====
def timeevolutionoperator(id,d_bare,modes_num,fields,t1,t2,U,info):
    id: atom_c_T instance
    d_bare: int
        Hilbert space dimension of the static system
    modes_num: 1D np.array([],dtype=np.int32)
        define the number of fundamental frequencies and modes
        the total number of modes is nm = np.sum(modes_num)
    fields: instance of openmmf.mode_c_T*nm
        equivalent to the FORTRAN data derived type FIELDS
    t1: np.double
        instant of time to evaluate the time evolution operator
    t2: np.double
        instant of time to evaluate the time evolution operator
    U: 1D np.array([],dtype=np.complex)
        d_bare x d_bare Matrix of the time-evolution operator
    info: int
        Error/success flag

=====
# // DEFINITION OF DRESSED BASIS WITH ALL FIELDS
=====
def dressedbasis(h_floquet_size,id,modes_num,fields,U_FD,e_dressed,info):
    h_floquet_size: int
        size of the multimode Floquet matrix
    id: atom_c_T instance
    modes_num: 1D np.array([],dtype=np.int32)
        define the number of fundamental frequencies and modes
        the total number of modes is nm = np.sum(modes_num)
    fields: instance of openmmf.mode_c_T*nm
        equivalent to the FORTRAN data derived type FIELDS
    U_FD: 1D np.array([],dtype=np.complex)
        d_bare x d_bare Matrix of the time-average state occupationrs
    e_dressed: 1D np.array([],dtype=np.complex)
        array containing the eigenvalues of the multimode Floquet
        Hamiltonian
    info: int
        Error/success flag

def dressedbasis_sp(h_floquet_size,id,modes_num,fields,U_FD,e_dressed,info):
    h_floquet_size: int
        size of the multimode Floquet matrix
    id: atom_c_T instance
    modes_num: 1D np.array([],dtype=np.int32)
        define the number of fundamental frequencies and modes
        the total number of modes is nm = np.sum(modes_num)
    fields: instance of openmmf.mode_c_T*nm
        equivalent to the FORTRAN data derived type FIELDS
    U_FD: 1D np.array([],dtype=np.complex)
        d_bare x d_bare Matrix of the time-average state occupationrs
    e_dressed: 1D np.array([],dtype=np.complex)
        array containing the eigenvalues of the multimode Floquet
        Hamiltonian
    info: int
        Error/success flag

=====

```

```

# // DEFINITION OF DRESSED BASIS WITH A SUBSET OF THE FIELDS
#=====
def dressedbasis_subset(id,dressingfields_indices,modes_num,fields,U_FD,e_dressed,info):
    id: atom_c_T instance
    dressingfields_indices: 1D np.array([],dtype=np.int32)
        define the index of dressing fields
    modes_num: 1D np.array([],dtype=np.int32)
        define the number of fundamental frequencies and modes
        the total number of modes is nm = np.sum(modes_num)
    fields: instance of openmmf.mode_c_T*nm
        equivalent to the FORTRAN data derived type FIELDS
    U_FD: 1D np.array([],dtype=np.complex)
        h_floquet_size x h_floquet_size Matrix of eigenvectors
    e_dressed: 1D np.array([],dtype=np.complex)
        array containing the eigenvalues of the multimode Floquet
        Hamiltonian
    info: int
        Error/success flag

def dressedbasis_subset_sp(id,dressingfields_indices,modes_num,fields,U_FD,e_dressed,info):
    id: atom_c_T instance
    dressingfields_indices: 1D np.array([],dtype=np.int32)
        define the index of dressing fields
    modes_num: 1D np.array([],dtype=np.int32)
        define the number of fundamental frequencies and modes
        the total number of modes is nm = np.sum(modes_num)
    fields: instance of openmmf.mode_c_T*nm
        equivalent to the FORTRAN data derived type FIELDS
    U_FD: 1D np.array([],dtype=np.complex)
        h_floquet_size x h_floquet_size Matrix of eigenvectors
    e_dressed: 1D np.array([],dtype=np.complex)
        array containing the eigenvalues of the multimode Floquet
        Hamiltonian
    info: int
        Error/success flag

#=====
# EVALUATE THE FOURIER COMPONENTS OF THE MICROMOTION OPERATOR USING THE DRESSING FIELDS
#=====
def micromotionfourierdressedbasis(id,dressingfields_indices,modes_num,fields,
    U_FD,e_dressed,info):
    id: atom_c_T instance
    dressingfields_indices: 1D np.array([],dtype=np.int32)
        define the index of dressing fields
    modes_num: 1D np.array([],dtype=np.int32)
        define the number of fundamental frequencies and modes
        the total number of modes is nm = np.sum(modes_num)
    fields: instance of openmmf.mode_c_T*nm
        equivalent to the FORTRAN data derived type FIELDS
    U_FD: 1D np.array([],dtype=np.complex)

    e_dressed: 1D np.array([],dtype=np.complex)
        array containing the eigenvalues of the multimode Floquet
        Hamiltonian
    info: int
        Error/success flag

```

```

#=====

```

```

# EVALUATE THE MICROMOTION OPERATOR USING THE DRESSING FIELDS
#=====
def micromotiondressedbasis(id,modes_num,dressingfields_indices,fields,t1,U,info):
    id: atom_c_T instance
    modes_num: 1D np.array([],dtype=np.int32)
        define the number of fundamental frequencies and modes
        the total number of modes is nm = np.sum(modes_num)
    dressingfields_indices: 1D np.array([],dtype=np.int32)
        define the index of dressing fields
    fields: instance of openmmf.mode_c_T*nm
        equivalent to the FORTRAN data derived type FIELDS
    t1: np.double
        instant of time to evaluate the micromotion operator
    U : 1D np.array([],dtype=np.complex)

    info: int
        Error/success flag

#=====
# DEALLOCATE ALL MEMORY ARRAYS
#=====
def deallocateall(id):
    id: atom_c_T instance

```