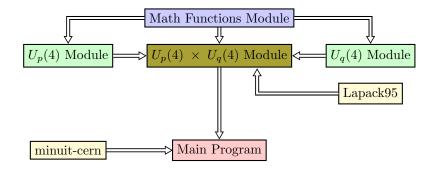
$$U_p(4) \times U_q(4)$$

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0.1 Math Functions Module \rightarrow MOD_matfun.f90

• Functions:

$$- \ {\tt p_symbol}({\tt a}, {\tt b}) = \ (a)_s \ = \ a(a+1)...(a+s-1)$$

- factorial(n) =
$$n!$$

- delta_function
$$(a,b,c) = \Delta(abc)$$

- wigner_6j
$$(j_1, j_2, j_3, l_1, l_2, l_3) = \begin{cases} j_1 & j_2 & j_3 \\ l_1 & l_2 & l_3 \end{cases}$$
 1

$$-\ \mathtt{wigner_9j}(j_1,j_2,j_3,l_1,l_2,l_3,k_1,k_2,k_3)\ =\ \begin{cases} j_1 & j_2 & j_3 \\ l_1 & l_2 & l_3 \\ k_1 & k_2 & k_3 \end{cases} \}^2$$

$0.2 \quad U_p(4) \text{ Module} \rightarrow \texttt{MOD_Up4.f90}$

Hamiltonian:

$$\hat{H}_{U_{p}(4)} = \beta \,\, \hat{\mathcal{C}}_{2} \left[so_{p}(4) \right] + \gamma \,\, \hat{\mathcal{C}}_{2} \left[so_{p}(3) \right] + \gamma_{2} \,\, \left[\hat{\mathcal{C}}_{2} \left[so_{p}(3) \right] \right]^{2} + \kappa \,\, \hat{\mathcal{C}}_{2} \left[so_{p}(4) \right] \hat{\mathcal{C}}_{2} \left[so_{p}(3) \right] \tag{1}$$

• Global definitions:

- Npval: U(4) Totally symmetric representation.

• Functions:

- Function: RME_Casimir_SOp4

- Function: RME_Casimir_SOp3

- Function: RME_Qp2

- Function: RME_np

 2 Integers only

 $^{^{1}}$ Definition of the book $\it Nuclear~Shell~Theory$ of Amos de-Shalit and Igal Talmi

0.3 $U_q(4)$ **Module** \rightarrow MOD_Uq4.f90

Hamiltonian:

$$\hat{H}_{U_q(4)} = a \,\hat{\mathcal{C}}_1 \left[u_q(3) \right] + b \,\hat{\mathcal{C}}_2 \left[u_q(3) \right] + c \,\hat{\mathcal{C}}_2 \left[so_q(3) \right] + d \,\hat{\mathcal{C}}_2 \left[so_q(4) \right] \tag{2}$$

- Global definitions:
 - Ngval: U(4) Totally symmetric representation.
- Functions:
 - Function: RME_Casimir_Uq3
 - Function: RME_Casimir_SOq3
 - Function: RME_Casimir_SOq4
 - Function: RME_Qq2
 - Function: RME_Dq_prima

$0.4 \quad U_p(4) \times U_q(4) \text{ Module} \rightarrow \text{MOD_Up_x_Uq.f90}$

- Global definitions:
 - basis_para(1:5, no of para-states): Integers. Para-states
 - basis_ortho(1:5, no of ortho-states): Integers. Ortho-states
 - $dim_para(1:\Lambda_{max})$: Integers. Para-dim blocks
 - dim_ortho(1: Λ_{max}): Integers. Ortho-dim blocks
 - $ijk_{para}(1:\Lambda_{max})$: Integers. Pseudo-pointer
 - ijk_ortho(1: Λ_{max}): Integers. Pseudo-pointer
 - lambda_max: Integer. Maximum value of Λ
 - Type exp_point: exp_point%ist, exp_point%i_pos,exp_point%fst,exp_point%f_pos, exp_point%energy and exp_point%intensity
 - total_exp: Integer. Number of experimental data.
 - exp_data: Type(exp_point)
 - Type matrix: $matrix(\Lambda)\%[SYM]$
 - $Ham(1:\Lambda_{max}): Type(matrix)$. Hamiltonian para/ortho matrices
 - SOq4(1: Λ_{max}): Type(matrix). $\hat{\mathcal{C}}_2[SO_q(4)]$ para/ortho matrices
 - QpQq(1: Λ_{\max}): Type(matrix). $\left[\hat{Q}_p^{(2)} \times \hat{Q}_q^{(2)}\right]^{(0)}$ para/ortho matrices
 - QpQqW(1: Λ_{max}): Type(matrix). $\left[\hat{Q}_p^{(2)} \times \hat{Q}_q^{(2)}\right]^{(0)} \hat{\mathcal{C}}_2 \left[SO_p(4)\right] + c.c.$ para/ortho matrices

• Functions:

- Subroutine: dimension_po

- Subroutine: build_basis_po

- Subroutine: initialize_position_index

Function: RME_Qp_x_Qq_0Function: RME_Ip_x_S0q4

- Function: RME_QpQqSOp4

- Subroutine: build_Up_x_Uq_matrix

Function: pretty_braketSubroutine: read_expdat

Function: exp_linesSubroutine: build_hamFunction: find_pos

Subroutine: eigensystemFunction: assig_state

Tunction, abbig_r

Function: chi2Function: FCN

- Function: RME_np_x_Dq_1

- Function: RME_Qp2_x_Dq_1

0.4.1 Building the basis

We have a block-diagonalizable system dividing the problem in para (mod(J, 2) = 0) and ortho (mod(J, 2) = 1) cases, and separating by different Λ . The states will be stored in the same array sorted by Λ . The dimension of each block will be saved in dim_[SYM] $(\Lambda = 0)$, ..., dim_[SYM] $(\Lambda = \Lambda_{\text{max}})$.

$$\begin{bmatrix} \left| \psi_1^{\Lambda=0} \right\rangle \\ \vdots \\ \left| \psi_{\text{dim_[SYM]}(\Lambda=0)}^{\Lambda=0} \right\rangle \end{bmatrix} \right\} \rightarrow 1: \text{dim_[SYM]}(\Lambda=0)$$

$$\vdots$$

$$\begin{bmatrix} \left| \psi_1^{\Lambda=\Lambda_{\text{max}}} \right\rangle \\ \vdots \\ \left| \psi_{\text{dim_[SYM]}(\Lambda=\Lambda_{\text{max}})}^{\Lambda=\Lambda_{\text{max}}} \right\rangle \end{bmatrix} \right\} \rightarrow 1: \text{dim_[SYM]}(\Lambda=\Lambda_{\text{max}})$$

$$(3)$$

Therefore, the basis array is as follow:

$$\mathtt{basis_[SYM]} = \left[\left| \psi_1^{\Lambda=0} \right\rangle, ..., \psi_{\mathtt{dim_[SYM]}(\Lambda=0)}^{\Lambda=0}, ..., \left| \psi_1^{\Lambda=\Lambda_{\max}} \right\rangle, ..., \psi_{\mathtt{dim_[SYM]}(\Lambda=\Lambda_{\max})}^{\Lambda=\Lambda_{\max}} \right] \tag{4}$$

Each element of the the basis must cotain information about quantum numbers:

$$|\psi_i^{\Lambda}\rangle = (w\ J;\ n\ L;\ \Lambda \in [|J-L|, J+L])$$
 (5)

Dimension: dimension_po

Fortran 90 subroutine. Inputs:

- Npval. Integer
- Nqval. Integer
- Λ_{max} . Integer

Outputs:

- dim_para. Integer, dimension $(0:\Lambda_{max})$
- dim_ortho. Integer, dimension $(0:\Lambda_{max})$

Basis: build_basis_po

Inputs:

- Npval. Integer
- Nqval. Integer
- $\Lambda_{\rm max}$. Integer
- total_para. Integer, total dimension of para basis
- total_ortho. Integer, total dimension of ortho basis
- dim_para. Integer, dimension($0:\Lambda_{max}$)
- dim_ortho. Integer, dimension $(0:\Lambda_{max})$

Outputs:

- basis_para. Integer, dimension $\left(1:5,\ 1:\sum_{\Lambda=0}^{\Lambda_{\max}} \mathtt{dim_para}(\Lambda)\right)$
- basis_ortho. Integer, dimension $\left(1:5,\ 1:\sum\limits_{\Lambda=0}^{\Lambda_{\max}} \mathtt{dim_ortho}(\Lambda)\right)$
 - basis_[SYM](1,:) $\longrightarrow w$
 - basis_[SYM](2,:) $\longrightarrow J$
 - basis_[SYM](3,:) $\longrightarrow n$
 - basis_[SYM](4,:) $\longrightarrow L$
 - $\text{ basis}_{\text{SYM}}(5,:) \longrightarrow \Lambda$

Para-states

$$\begin{split} \text{LOOP: } J &= 0, 2, ..., N_p - \operatorname{mod}(N_p, 2) \\ \text{LOOP: } L &= 0, 1, ..., N_q \\ \text{CONDITIONAL: } |J - L| \leq \Lambda_{max} \text{ to continue, else go to next } L \\ \text{LOOP: } \Lambda &= |J - L|, |J - L| + 1, ..., \min(\Lambda_{max}, J + L) \\ \text{LOOP: } w &= N_p - \operatorname{mod}(N_p, 2), N_p - \operatorname{mod}(N_p, 2) - 2, ..., J \\ \text{LOOP: } n &= L, L + 2, ..., Nq \\ \text{basis_para}\left(1, \operatorname{dim_para}\left(\Lambda\right)\right) &= w \\ \text{basis_para}\left(2, \operatorname{dim_para}\left(\Lambda\right)\right) &= J \\ \text{basis_para}\left(3, \operatorname{dim_para}\left(\Lambda\right)\right) &= L \\ \text{basis_para}\left(4, \operatorname{dim_para}\left(\Lambda\right)\right) &= L \\ \text{basis_para}\left(5, \operatorname{dim_para}\left(\Lambda\right)\right) &= \Lambda \end{split}$$

Ortho-states

$$\begin{split} \text{LOOP: } J &= 1, 3, ..., N_p - (1 - \operatorname{mod}(N_p, 2)) \\ \text{LOOP: } L &= 0, 1, ..., N_q \\ \text{CONDITIONAL: } |J - L| \leq \Lambda_{max} \text{ to continue, else go to next } L \\ \text{LOOP: } \Lambda &= |J - L|, |J - L| + 1, ..., \min(\Lambda_{max}, J + L) \\ \text{LOOP: } w &= N_p - (1 - \operatorname{mod}(N_p, 2)), N_p - (1 - \operatorname{mod}(N_p, 2)) - 2, ..., J \\ \text{LOOP: } n &= L, L + 2, ..., Nq \\ \text{basis_ortho} (1, \operatorname{dim_ortho}(\Lambda)) &= w \\ \text{basis_ortho} (2, \operatorname{dim_ortho}(\Lambda)) &= J \\ \text{basis_ortho} (3, \operatorname{dim_ortho}(\Lambda)) &= L \\ \text{basis_ortho} (5, \operatorname{dim_ortho}(\Lambda)) &= L \\ \text{basis_ortho} (5, \operatorname{dim_ortho}(\Lambda)) &= \Lambda \end{split}$$

${\bf Pseudo-pointer:\ initialize_position_index}$

```
subroutine initialize_position_index(ijk,partial_dim,lambda_max)
!
! Inputs:
! o) partial_dim: Integer array dimension 0:lambda_max
! o) lambda_max
!
! Output:
! o) ijk: Integer array dimension 0:lambda_max
!
! This functions initializes the initial integer "pointer" ijk(0:lambda_max),
! so that
! ijk(0) = 1
! ijk(1) = position where lambda=1 block starts
! ...
! ijk(lambda_max) = position where lambda=lambda_max block stats
!
! This function is going to be of vital importance during the program's development
!
```

Building all matrices: build_Up_x_Uq_matrix

This subroutine builds the matrices for the operators of $U_p(4) \times U_q(4)$ using

The function RME_fun must deppend on $(\omega_1, J_1, n_1, L_1, \Lambda_1, \omega_2, J_2, n_2, L_2, \Lambda_2)$.

! All position corresponding to lam1 /= lam2 will be ZERO!

Braket notation output: pretty_braket

```
Useless gadget ... but very nice.

function pretty_braket(w,j,n,l,lam,bk,Np,Nq)

!
! INPUTs:
! o) Np(opt), Nq(opt), w, j, n, l, lam: Quantum numbers
! o) bk: one character = b (bra) or k (ket)
!
! OUTPUT:
! o) pretty_braket: character type
```