Program Reference

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Interface

C routine

```
dim = CINTgto_cart(bas_id, bas);
dim = CINTgto_spheric(bas_id, bas);
dim = CINTgto_spinor(bas_id, bas);
f1e(buf, shls, atm, natm, bas, nbas, env);
f2e(buf, shls, atm, natm, bas, nbas, env, opt);
f2e_optimizer(&opt, atm, natm, bas, nbas, env);
CINTdel_optimizer(&opt);
```

- buf: column-major double precision array.
  - for 1e integrals of shells (i,j), data are stored as `[i1j1 i2j1 ... ]`
  - for 2e integrals of shells (i,j|k,l), data are stored as
    `[i1j1k1l1 i2j1k1l1 ... i1j2k1l1 ... i1j1k2l1 ... ]`

    - complex data are stored as two double elements, first is real, followed by imaginary, e.g. [Re Im Re Im …]
- shls: 0-based basis/shell indices.
    - int[2] for 1e integrals
    - int[4] for 2e integrals
- atm: int[natm*6], list of atoms. For ith atom, the 6 slots of atm[i] are
    - `atm[i*6+0]` nuclear charge of atom i
    - `atm[i*6+1]` env offset to save coordinates (`env[atm[i*6+1]]`, `env[atm[i*6+1]+1]`, `env[atm[i*6+1]+2]`) are (x,y,z)
    - `atm[i*6+2]` nuclear model of atom i, = 2 indicates gaussian nuclear model $\rho(r) = Z(\frac{\zeta}{\pi})^{3/2}\exp(-\zeta r^2)$
    - `atm[i*6+3]` env offset to save the nuclear charge distribution parameter ζ
    - `atm[i*6+4]` unused
    - `atm[i*6+5]` unused
- natm: int, number of atoms, natm has no effect **except nuclear attraction** integrals
- bas: int[nbas*8], list of basis. For ith basis, the 8 slots of bas[i] are
    - `bas[i*8+0]` 0-based index of corresponding atom
    - `bas[i*8+1]` angular momentum
    - `bas[i*8+2]` number of primitive GTO in basis i
    - `bas[i*8+3]` number of contracted GTO in basis i
    - `bas[i*8+4]` kappa for spinor GTO.
      $<$ 0 the basis ~ j = l + 1/2.
      $>$ 0 the basis ~ j = l - 1/2.
      = 0 the basis includes both j = l + 1/2 and j = l - 1/2
    - `bas[i*8+5]` env offset to save exponents of primitive GTOs. e.g. 10 exponents `env[bas[i*8+5]]` … `env[bas[i*8+5]+9]`
    - `bas[i*8+6]` env offset to save column-major contraction coefficients. e.g. 10 primitive -> 5 contraction needs a 10 × 5 array

```
env[bas[i*8+6]   ] | env[bas[i*8+6]+10] |     | env[bas[i*8+6]+40]
env[bas[i*8+6]+1]  | env[bas[i*8+6]+11] |     | env[bas[i*8+6]+41]
 .                 |  .                 | ... |  .
 .                 |  .                 |     |  .
env[bas[i*8+6]+9]  | env[bas[i*8+6]+19] |     | env[bas[i*8+6]+49]
```

```
- ‘bas[i*8+7]‘ unused
```

- nbas: int, number of bases, nbas has no effect, can be set to 0
- env: double[], save the value of coordinates, exponents, contraction coefficients
- struct CINTOpt *opt: so called “optimizer”, it needs to be intialized

  CINTOpt *opt = NULL; intname_optimizer(&opt, atm, natm, bas, nbas, env);

every integral type has its own optimizer with the suffix _optimizer in its name, e.g. the optimizer for cint2e_sph is cint2e_sph_opimizer. “optimizer” is an optional argument for the integrals. It can roughly speed the integration by 10% without affecting the value of integrals. If no optimizer is wanted, set it to NULL.

optimizer needs to be released after using.

CINTdel_optimizer(&opt);

- if the return value equals 0, every element of the integral is 0
- short example

```
#include "cint.h"
...
CINTOpt *opt = NULL;
cint2e_sph_optimizer(&opt, atm, natm, bas, nbas, env);
for (i = 0; i < nbas; i++) {
        shls[0] = i;
        di = CINTcgto_spheric(i, bas);
        ...
        for (l = 0; l < nbas; l++) {
                shls[3] = l;
                dl = CINTcgto_spheric(l, bas);
                buf = malloc(sizeof(double) * di * dj * dk * dl);
                cint2e_cart(buf, shls, atm, natm, bas, nbas, env, opt);
                free(buf);
        }
}
CINTdel_optimizer(&opt);
```

Fortran routine

```
dim = CINTgto_cart(bas_id, bas)
dim = CINTgto_spheric(bas_id, bas)
dim = CINTgto_spinor(bas_id, bas)
call f1e(buf, shls, atm, natm, bas, nbas, env)
call f2e(buf, shls, atm, natm, bas, nbas, env, opt)
call f2e_optimizer(opt, atm, natm, bas, nbas, env)
call CINTdel_optimizer(opt)
```

- atm and bas are 2D integer array
  - atm(1:6,i) is the (charge, offset_coord, nuclear_model, unused, unused, unused) of the ith atom
  - bas(1:8,i) is the (atom_index, angular, num_primitive_GTO, num_contract_GTO, kappa, offset_exponent, offset_coeff, unused) of the ith basis

- parameters are the same to the C function. Note that those offsets atm(2,i) bas(6,i) bas(7,i) are 0-based.
- buf is 2D/4D double precision/double complex array
- opt: an integer(8) to hold the address of so called "optimizer", it needs to be intialized by

  integer(8) opt call f2e_optimizer(opt, atm, natm, bas, nbas, env)

  The optimizier can be banned by setting the "optimizier" to 0_8

  call f2e(buf, atm, natm, bas, nbas, env, 0_8)

  To release optimizer, execute

  call CINTdel_optimizer(opt);
- short example

```
...
integer,external CINTcgto_spheric
integer(8) opt
call cint2e_sph_optimizer(opt, atm, natm, bas, nbas, env)
do i = 1, nbas
  shls(1) = i - 1
  di = CINTcgto_spheric(i-1, bas)
  ...
  do l = 1, nbas
    shls(4) = l - 1
    dl = CINTcgto_spheric(l-1, bas)
    allocate(buf(di,dj,dk,dl))
    call cint2e_sph(buf, shls, atm, natm, bas, nbas, env, opt)
    deallocate(buf)
  end do
end do
call CINTdel_optimizer(opt)
```

Supported angular momentum

$l_{max} = 6$

Data ordering

- for Cartesian GTO, the output data in buf are sorted as

s shell	p shell	d shell	...
...	...	...	
s	p x	d xx	
s	p y	d xy	
...	p z	d xz	
	p x	d yy	
	p y	d yz	
	p z	d zz	
	...	...	

- for real spheric GTO, the output data in buf are sorted as

s shell	p shell	d shell	f shell	...
...	...	...	...	
s	p x	d xy	f $y(3x^2-y^2)$	
s	p y	d yz	f xyz	
...	p z	d z^2	f yz^2	
	p x	d xz	f z^3	
	p y	d x^2-y^2	f xz^2	
	p z	...	f $z(x^2-y^2)$	
	...		f $x(x^2-3y^2)$	
			...	

- for spinor GTO, the output data in buf correspond to

...	kappa=0,p shell	kappa=1,p shell	kappa=0,d shell	...
	...	...	...	
	$p_{1/2}(-1/2)$	$p_{1/2}(-1/2)$	$d_{3/2}(-3/2)$	
	$p_{1/2}(1/2)$	$p_{1/2}(1/2)$	$d_{3/2}(-1/2)$	
	$p_{3/2}(-3/2)$	$p_{1/2}(-1/2)$	$d_{3/2}(1/2)$	
	$p_{3/2}(-1/2)$	$p_{1/2}(1/2)$	$d_{3/2}(3/2)$	
	$p_{3/2}(1/2)$	$p_{1/2}(-1/2)$	$d_{5/2}(-5/2)$	
	$p_{3/2}(3/2)$	$p_{1/2}(1/2)$	$d_{5/2}(-3/2)$	
	$p_{1/2}(-1/2)$	...	$d_{5/2}(-1/2)$	
	$p_{1/2}(1/2)$		$d_{3/2}(-3/2)$	
	$p_{3/2}(-3/2)$		$d_{3/2}(-1/2)$	
	$p_{3/2}(-1/2)$		...	
	...			

Tensor

Integrals like Gradients have more than one components. The output array is ordered in Fortran-contiguous. The tensors are ordered as

- 3-component tensor
  - X `buf(:,0)`
  - Y `buf(:,1)`
  - Z `buf(:,2)`
- 9-component tensor

- XX `buf(:,0)`
- XY `buf(:,1)`
- XZ `buf(:,2)`
- YX `buf(:,3)`
- YY `buf(:,4)`
- YZ `buf(:,5)`
- ZX `buf(:,6)`
- ZY `buf(:,7)`
- ZZ `buf(:,8)`

Built-in function list

- Cartesian GTO integrals
  - `CINTcgto_cart(int shell_id, int bas[])`: Number of cartesian functions of the given shell
  - `cint1e_ovlp_cart`
    $$\langle i|j\rangle$$
  - `cint1e_nuc_cart`
    $$\langle i|V_{nuc}|j\rangle$$
  - `cint1e_kin_cart`
    $$.5\langle i|\vec{p}\cdot\vec{p}j\rangle$$
  - `cint1e_ia01p_cart`
    $$\langle i|\frac{\vec{r}}{r^3}|\times\vec{\nabla}j\rangle$$
  - `cint1e_irixp_cart`
    $$\langle i|(\vec{r}-\vec{R}_i)\times\vec{\nabla}j\rangle$$
  - `cint1e_ircxp_cart`
    $$\langle i|(\vec{r}-\vec{R}_o)\times\vec{\nabla}j\rangle$$
  - `cint1e_iking_cart`
    $$0.5i\langle\vec{p}\cdot\vec{p}i|U_g j\rangle$$
  - `cint1e_iovlpg_cart`
    $$i\langle i|U_g j\rangle$$
  - `cint1e_inucg_cart`
    $$i\langle i|V_{nuc}|U_g j\rangle$$
  - `cint1e_ipovlp_cart`
    $$\langle\vec{\nabla}i|j\rangle$$
  - `cint1e_ipkin_cart`
    $$0.5\langle\vec{\nabla}i|\vec{p}\cdot\vec{p}j\rangle$$
  - `cint1e_ipnuc_cart`
    $$\langle\vec{\nabla}i|V_{nuc}|j\rangle$$

- `cint1e_iprinv_cart`
$$\langle\vec{\nabla}i|r^{-1}|j\rangle$$
- `cint1e_rinv_cart`
$$\langle i|r^{-1}|j\rangle$$
- `cint2e_cart`
$$(ij|kl)$$
- `cint2e_ig1_cart`
$$i(iU_gj|kl)$$
- `cint2e_ip1_cart`
$$(\vec{\nabla}ij|kl)$$

• Spheric GTO integrals
  - `CINTcgto_spheric(int shell_id, int bas[])`: Number of spheric functions of the given shell
  - `cint1e_ovlp_sph`
$$\langle i|j\rangle$$
  - `cint1e_nuc_sph`
$$\langle i|V_{nuc}|j\rangle$$
  - `cint1e_kin_sph`
$$0.5\langle i|\vec{p}\cdot pj\rangle$$
  - `cint1e_ia01p_sph`
$$\langle i|\frac{\vec{r}}{r^3}|\times\vec{\nabla}j\rangle$$
  - `cint1e_irixp_sph`
$$\langle i|(\vec{r}_c-\vec{R}_i)\times\vec{\nabla}j\rangle$$
  - `cint1e_ircxp_sph`
$$\langle i|(\vec{r}_c-\vec{R}_o)\times\vec{\nabla}j\rangle$$
  - `cint1e_iking_sph`
$$0.5i\langle\vec{p}\cdot\vec{p}i|U_gj\rangle$$
  - `cint1e_iovlpg_sph`
$$i\langle i|U_gj\rangle$$
  - `cint1e_inucg_sph`
$$i\langle i|V_{nuc}|U_gj\rangle$$
  - `cint1e_ipovlp_sph`
$$\langle\vec{\nabla}i|j\rangle$$
  - `cint1e_ipkin_sph`
$$0.5\langle\vec{\nabla}i|\vec{p}\cdot pj\rangle$$
  - `cint1e_ipnuc_sph`
$$\langle\vec{\nabla}i|V_{nuc}|j\rangle$$

  - `cint1e_iprinv_sph`
$$\langle \vec{\nabla} i | r^{-1} | j \rangle$$
  - `cint1e_rinv_sph`
$$\langle i | r^{-1} | j \rangle$$
  - `cint2e_sph`
$$(ij|kl)$$
  - `cint2e_ig1_sph`
$$i(iU_g j|kl)$$
  - `cint2e_ip1_sph`
$$(\vec{\nabla} ij|kl)$$

- Spinor GTO integrals
  - `CINTcgto_spinor(int shell_id, int bas[])`: Number of spinor functions of the given shell
  - `cint1e_ovlp`
$$\langle i | j \rangle$$
  - `cint1e_nuc`
$$\langle i | V_{nuc} | j \rangle$$
  - `cint1e_nucg`
$$\langle i | V_{nuc} | U_g j \rangle$$
  - `cint1e_srsr`
$$\langle \vec{\sigma} \cdot \vec{r} i | \vec{\sigma} \cdot \vec{r} j \rangle$$
  - `cint1e_sr`
$$\langle \vec{\sigma} \cdot \vec{r} i | j \rangle$$
  - `cint1e_srsp`
$$\langle \vec{\sigma} \cdot \vec{r} i | \vec{\sigma} \cdot \vec{p} j \rangle$$
  - `cint1e_spsp`
$$\langle \vec{\sigma} \cdot \vec{p} i | \vec{\sigma} \cdot \vec{p} j \rangle$$
  - `cint1e_sp`
$$\langle \vec{\sigma} \cdot \vec{p} i | j \rangle$$
  - `cint1e_spspsp`
$$\langle \vec{\sigma} \cdot \vec{p} i | \vec{\sigma} \cdot \vec{p} \vec{\sigma} \cdot \vec{p} j \rangle$$
  - `cint1e_spnuc`
$$\langle \vec{\sigma} \cdot \vec{p} i | V_{nuc} | j \rangle$$
  - `cint1e_spnucsp`
$$\langle \vec{\sigma} \cdot \vec{p} i | V_{nuc} | \vec{\sigma} \cdot \vec{p} j \rangle$$
  - `cint1e_srnucsr`
$$\langle \vec{\sigma} \cdot \vec{r} i | V_{nuc} | \vec{\sigma} \cdot \vec{r} j \rangle$$

- `cint1e_sa10sa01`

$$0.5\langle\vec{\sigma}\times\vec{r}_c i|\vec{\sigma}\times\frac{\vec{r}}{r^3}|j\rangle$$

- `cint1e_ovlpg`

$$\langle i|U_g j\rangle$$

- `cint1e_sa10sp`

$$0.5\langle\vec{r}_c\times\vec{\sigma}i|\vec{\sigma}\cdot\vec{p}j\rangle$$

- `cint1e_sa10nucsp`

$$0.5\langle\vec{r}_c\times\vec{\sigma}i|V_{nuc}|\vec{\sigma}\cdot\vec{p}j\rangle$$

- `cint1e_sa01sp`

$$\langle i|\frac{\vec{r}}{r^3}\times\vec{\sigma}|\vec{\sigma}\cdot\vec{p}j\rangle$$

- `cint1e_spgsp`

$$\langle U_g\vec{\sigma}\cdot\vec{p}i|\vec{\sigma}\cdot\vec{p}j\rangle$$

- `cint1e_spgnucsp`

$$\langle U_g\vec{\sigma}\cdot\vec{p}i|V_{nuc}|\vec{\sigma}\cdot\vec{p}j\rangle$$

- `cint1e_spgsa01`

$$\langle U_g\vec{\sigma}\cdot\vec{p}i|\frac{\vec{r}}{r^3}\times\vec{\sigma}|j\rangle$$

- `cint1e_ipovlp`

$$\langle\vec{\nabla}i|j\rangle$$

- `cint1e_ipkin`

$$0.5\langle\vec{\nabla}i|p\cdot pj\rangle$$

- `cint1e_ipnuc`

$$\langle\vec{\nabla}i|V_{nuc}|j\rangle$$

- `cint1e_iprinv`

$$\langle\vec{\nabla}i|r^{-1}|j\rangle$$

- `cint1e_ipspnucsp`

$$\langle\vec{\nabla}\vec{\sigma}\cdot\vec{p}i|V_{nuc}|\vec{\sigma}\cdot\vec{p}j\rangle$$

- `cint1e_ipsprinvsp`

$$\langle\vec{\nabla}\vec{\sigma}\cdot\vec{p}i|r^{-1}|\vec{\sigma}\cdot\vec{p}j\rangle$$

- `cint2e`

$$(ij|kl)$$

- `cint2e_spsp1`

$$(\vec{\sigma}\cdot\vec{p}i\vec{\sigma}\cdot\vec{p}j|kl)$$

- `cint2e_spsp1spsp2`

$$(\vec{\sigma}\cdot\vec{p}i\vec{\sigma}\cdot\vec{p}j|\vec{\sigma}\cdot\vec{p}k\vec{\sigma}\cdot\vec{p}l)$$

- `cint2e_srsr1`
$$(\vec{\sigma}\cdot\vec{r}i\vec{\sigma}\cdot\vec{r}j|kl)$$
- `cint2e_srsr1srsr2`
$$(\vec{\sigma}\cdot\vec{r}i\vec{\sigma}\cdot\vec{r}j|\vec{\sigma}\cdot\vec{r}k\vec{\sigma}\cdot\vec{r}l)$$
- `cint2e_sa10sp1`
$$0.5(\vec{r}_c\times\vec{\sigma}i\vec{\sigma}\cdot\vec{p}j|kl)$$
- `cint2e_sa10sp1spsp2`
$$0.5(\vec{r}_c\times\vec{\sigma}i\vec{\sigma}\cdot\vec{p}j|\vec{\sigma}\cdot\vec{p}k\vec{\sigma}\cdot\vec{p}l)$$
- `cint2e_g1`
$$(iU_gj|kl)$$
- `cint2e_spgsp1`
$$(\vec{\sigma}\cdot\vec{p}iU_g\vec{\sigma}\cdot\vec{p}j|kl)$$
- `cint2e_g1spsp2`
$$(iU_gj|\vec{\sigma}\cdot\vec{p}k\vec{\sigma}\cdot\vec{p}l)$$
- `cint2e_spgsp1spsp2`
$$(\vec{\sigma}\cdot\vec{p}iU_g\vec{\sigma}\cdot\vec{p}j|\vec{\sigma}\cdot\vec{p}k\vec{\sigma}\cdot\vec{p}l)$$
- `cint2e_ip1`
$$(\vec{\nabla}ij|kl)$$
- `cint2e_ipspsp1`
$$(\vec{\nabla}\vec{\sigma}\cdot\vec{p}i\vec{\sigma}\cdot\vec{p}j|kl)$$
- `cint2e_ip1spsp2`
$$(\vec{\nabla}ij|\vec{\sigma}\cdot\vec{p}k\vec{\sigma}\cdot\vec{p}l)$$
- `cint2e_ipspsp1spsp2`
$$(\vec{\nabla}\vec{\sigma}\cdot\vec{p}i\vec{\sigma}\cdot\vec{p}j|\vec{\sigma}\cdot\vec{p}k\vec{\sigma}\cdot\vec{p}l)$$
- `cint2e_ssp1ssp2`
$$(i\vec{\sigma}\vec{\sigma}\cdot\vec{p}j|k\vec{\sigma}\vec{\sigma}\cdot\vec{p}l)$$