

# 1 Total Energy for GFN2-xTB

$$\begin{aligned} E_{GFN2-xTB} &= E_{rep}^{(0)} + E_{disp}^{(0,1,2)} + E_{EHT}^{(1)} + E_{IES+IXC}^{(2)} + E_{AES+AXC}^{(2)} + E_{IES+IXC}^{(3)} \\ &= E_{rep} + E_{disp}^{D4'} + E_{EHT} + E_{\gamma} + E_{AES} + E_{AXC} + E_{\Gamma}^{GFN2} \end{aligned} \quad (1)$$

## 1.1 Repulsion Energy

$$E_{rep} = \frac{1}{2} \sum_{A,B} \frac{Z_A^{eff} Z_B^{eff}}{R_{AB}} e^{-\sqrt{a_A a_B} (R_{AB})^{(k_f)}} \quad (2)$$

$$k_f = \begin{cases} 1 & \text{if } A, B \in \{\text{H, He}\} \\ \frac{3}{2} & \text{otherwise} \end{cases} \quad (3)$$

$Z^{eff}$  and  $a$  are variables fitted for each element. A,B are the labels of atoms. Since we only have C and H in our systems we can simplify this quite a bit in code.  $R_{AB}$  is the distance between the A and B atoms.

## 1.2 Extended Hückel Theory Energy

$$E_{EHT} = \sum_{\mu\nu} P_{\mu\nu} H_{\mu\nu}^{EHT} \quad (4)$$

where  $\mu$  and  $\nu$  are AO indices,  $l$  and  $l'$  index shells. Both AO's are associated with an atom labeled A and B.

$$P_{\mu\nu} = P_{\mu\nu}^{(0)} + \delta P_{\mu\nu} \quad (5)$$

$$P_{\mu\nu}^{(0)} = \sum_j^{occ.MO} c_{\mu j} c_{j\nu} \quad (6)$$

$$\delta P_{\mu\nu} = ?? \quad \text{comes from the iteration, can be skipped for now} \quad (7)$$

$$\begin{aligned} H_{\mu\nu}^{EHT} &= \frac{1}{2} K_{AB}^{ll'} S_{\mu\nu} (H_{\mu\mu} + H_{\nu\nu}) \\ &\cdot X(EN_A, EN_B) \\ &\cdot \Pi(R_{AB}, l, l') \\ &\cdot Y(\zeta_l^A, \zeta_{l'}^B), \forall \mu \in l(A), \nu \in l'(B) \end{aligned} \quad (8)$$

$K_{AB}^{ll'}$  is a element and shell specific fitted constant however, in GFN2 it only depends on the shells.  $S_{\mu\nu} = \langle \phi_\mu | \phi_\nu \rangle$  is just the overlap of the orbitals. In GFN2  $H_{\kappa\kappa} = h_A^l - \delta h_{CN'_A}^l CN'_A$  where  $CN'_A$  is the modified GFN2-type Coordinate Number for the element of atom A.

$$\begin{aligned} CN'_A &= \sum_{B \neq A}^{N_{atoms}} (1 + e^{-10(4(R_{A,cov} + R_{B,cov})/3R_{AB} - 1)})^{-1} \\ &\times (1 + e^{-20(4(R_{A,cov} + R_{B,cov} + 2)/3R_{AB} - 1)})^{-1} \end{aligned} \quad (9)$$

$h_A^l$  and  $\delta h_{CN_A}^l$  are both fitted constants.  $EN_A$  is the electronegativity of the element of atom A.

$$X(EN_A, EN_B) = 1 + k_{EN} \Delta EN_{AB}^2 \quad (10)$$

$$k_{EN} = 0.02 \text{ in GFN2} \quad (11)$$

$$\Delta EN_{AB}^2 = (EN_A - EN_B)^2 \quad (12)$$

The electronegativity for C and H are 2.55 and 2.20 according to wikipedia. Thus here is a table for the combinations we will be working with:

A	B	$X(EN_A, EN_B)$
C	C	1
C	H	$1 + 0.02 \cdot (0.35^2)$
H	C	$1 + 0.02 \cdot (0.35^2)$
H	H	1

$$\Pi(R_{AB}, l, l') = \left( 1 + k_{A,l}^{\text{poly}} \left( \frac{R_{AB}}{R_{\text{cov},AB}} \right)^{\frac{1}{2}} \right) \left( 1 + k_{B,l'}^{\text{poly}} \left( \frac{R_{AB}}{R_{\text{cov},AB}} \right)^{\frac{1}{2}} \right) \quad (13)$$

$R_{\text{cov},AB}$  are the summed covalent radii ( $R_{\text{cov},A} + R_{\text{cov},B}$ ), the covalent radii is the second number in the table for each element,  $H_{\text{cov}} = 0.32$ ,  $C_{\text{cov}} = 0.75$  (can be found on wiki as 0.75).  $k_{A,l}^{\text{poly}}$  and  $k_{B,l'}^{\text{poly}}$  are element and shell specific constants.

$$Y(\zeta_l^A, \zeta_{l'}^B) = \left( \frac{2\sqrt{\zeta_l^A \zeta_{l'}^B}}{\zeta_l^A + \zeta_{l'}^B} \right)^{\frac{1}{2}} \quad (14)$$

Here,  $\zeta_l^A$  are the STO exponents of the GFN2-xTB AO basis. Slater Type Orbitals are defined as such:

$$\chi_{\zeta,n,l,m}(r, \theta, \varphi) = N Y_{l,m}(\theta, \varphi) r^{n-1} e^{-\zeta r} \quad (15)$$

N is a normalisation constant, Y are spherical harmonic funtions, n, l, m are the quantum numbers for the AO.  $r, \theta, \varphi$  are polar 3D coordinates.  $\zeta$  determines the radial extent of the STO, a large value gives rise to a function that is "tight" around the nucleus and a small value gives a more "diffuse" function. This  $\zeta$  is the one mentioned in the Y term of  $E_{EHT}$  and is a value fitted when constructing the basis set, thus it is given to us.

### 1.3 Isotropic electrostatic and XC energy

#### 1.3.1 Second order

$$E_\gamma = \frac{1}{2} \sum_{A,B} \sum_{l \in A} \sum_{l' \in B} q_l q_{l'} \gamma_{AB, ll'} \quad (16)$$

$$\gamma_{AB, ll'} = \frac{1}{\sqrt{R_{AB}^2 + \eta_{AB, ll'}^{-2}}} \quad (17)$$

$$\eta_{AB, ll'} = \frac{1}{2} \left[ \eta_A (1 + k_A^l) + \eta_B (1 + k_B^{l'}) \right] \quad (18)$$

$q_l$  is a partial muliken charge.  $\eta_A$  and  $\eta_B$  are element-specific fit parameters, while  $k_A^l$  and  $k_B^{l'}$  are element-specific scaling factors for the individual shells ( $k_A^l = 0$  when  $l = 0$ ).

$$GAP_A = \sum_{l \in A} q_l \quad (19)$$

$$q_l = \sum_{l' \in B} P_{ll'} S_{ll'} = GOP_l \quad (20)$$

#### 1.3.2 Third order

$$E_\Gamma^{GFN2} = \frac{1}{3} \sum_A \sum_{l \in A} (q_l)^3 K_l^\Gamma \Gamma_A \quad (21)$$

$K_l^\Gamma$  is a shell specific constant common for all elements and  $\Gamma_A$  is an element specific constant.

### 1.4 Anisotropic electrostatic energy

$$\begin{aligned} E_{AES} &= E_{q\mu} + E_{q\Theta} + E_{\mu\mu} \\ &= \frac{1}{2} \sum_{A,B} \{ f_3(R_{AB}) [q_A (\boldsymbol{\mu}_B^T \mathbf{R}_{BA}) + q_B (\boldsymbol{\mu}_A^T \mathbf{R}_{AB})] \\ &\quad + f_5(R_{AB}) [q_A \mathbf{R}_{AB}^T \boldsymbol{\Theta}_B \mathbf{R}_{AB} + q_B \mathbf{R}_{AB}^T \boldsymbol{\Theta}_A \mathbf{R}_{AB} \\ &\quad - 3(\boldsymbol{\mu}_A^T \mathbf{R}_{AB})(\boldsymbol{\mu}_B^T \mathbf{R}_{AB}) + (\boldsymbol{\mu}_A^T \boldsymbol{\mu}_B) R_{AB}^2] \} \end{aligned} \quad (22)$$

$\boldsymbol{\mu}_A$  is the cumulative atomic dipole moment of atom A and  $\boldsymbol{\Theta}_A$  is the corresponding traceless quadrupole moment. The curly braces and brackets are used in the same way as normal parenthesis for showing order of operations.  $q_A$  is

the atomic charge of atom A.

$$\Theta_A^{\alpha\beta} = \frac{3}{2}\theta_A^{\alpha\beta} - \frac{\delta_{\alpha\beta}}{2}(\theta_A^{xx} + \theta_A^{yy} + \theta_A^{zz}) \quad (23)$$

$$\theta_A^{\alpha\beta} = \sum_{l' \in A} \sum_l P_l \left( \alpha_A D_{ll'}^\beta + \beta_A D_{ll'}^\alpha - \alpha_A \beta_A S_{ll'} - Q_{ll'}^{\alpha\beta} \right) \quad (24)$$

$$q_A = Z_A - GAP_A \quad (25)$$

$$\mu_A^\alpha = \sum_{l' \in A} \sum_l P_{l'l} (\alpha_A S_{l'l} - D_{l'l}^\alpha) \quad (26)$$

$$D_{ll'}^\alpha = \langle \phi_l | \alpha_i | \phi_{l'} \rangle \quad (27)$$

$$Q_{ll'}^{\alpha\beta} = \langle \phi_l | \alpha_i \beta_i | \phi_{l'} \rangle \quad (28)$$

$\alpha$  and  $\beta$  are Cartesian components labled  $(x, y, z)^T$  with atom A being centered in  $v_A = (x_i, y_i, z_i)^T$  where i is a form of pointer/label dereferencing.  $\delta_{\alpha\beta}$  is just the delta function, i.e is 1 if  $\alpha$  and  $\beta$  are the same label and 0 otherwise, this serves to include the term only for the diagonal.

$$\Theta_A = \begin{pmatrix} \Theta_A^{xx} & \Theta_A^{xy} & \Theta_A^{xz} \\ \Theta_A^{yx} & \Theta_A^{yy} & \Theta_A^{yz} \\ \Theta_A^{zx} & \Theta_A^{zy} & \Theta_A^{zz} \end{pmatrix} \quad (29)$$

$$\mu_A = \begin{pmatrix} \mu_A^x \\ \mu_A^y \\ \mu_A^z \end{pmatrix} \quad (30)$$

$$R_{AB} = v_A - v_B \quad (31)$$

$$f_n(R_{AB}) = \frac{f_{damp}(a_n, R_{AB})}{R_{AB}^n} = \frac{1}{R_{AB}^n} \frac{1}{1 + 6 \left( \frac{R_0^{AB}}{R_{AB}} \right)^{a_n}} \quad (32)$$

$$R_0^{AB} = 0.5(R_0^{A'} + R_0^{B'}) \quad (33)$$

$$R_0^{A'} = \begin{cases} R_0^A + \frac{R_{max} - R_0^A}{1 + \exp[-4(CN_A' - N_{val} - \Delta_{val})]} & \text{if } N_{val} \text{ is given} \\ 5.0 \text{ bohrs} & \text{otherwise} \end{cases} \quad (34)$$

$$R_{max} = 5.0 \text{ bohrs} \quad (35)$$

$$\Delta_{val} = 1.2 \quad (36)$$

$R_0^A$  is a fitted value for 12 elements and 5.0 for the rest.  $a_n$  are adjusted global parameters.

## 1.5 Anisotropic XC energy

$$E_{AXC} = \sum_A (f_{XC}^{\mu_A} |\mu_A|^2 + f_{XC}^{\Theta_A} \|\Theta_A\|^2) \quad (37)$$

Where  $f_{XC}^{\mu_A}$  and  $f_{XC}^{\Theta_A}$  are fitted values. What norms are these?

## 1.6 Dispersion Energy

$$\begin{aligned}
E_{disp}^{D4'} = & - \sum_{A>B} \sum_{n=6,8} s_n \frac{C_n^{AB}(q_A, CN_{cov}^A, q_B, CN_{cov}^B)}{R_{AB}^n} f_{damp,BJ}^{(n)}(R_{AB}) \\
& - s_9 \sum_{A>B>C} \frac{(3\cos(\theta_{ABC})\cos(\theta_{BCA})\cos(\theta_{CAB}) + 1)C_9^{ABC}(CN_{cov}^A, CN_{cov}^B, CN_{cov}^C)}{(R_{AB}R_{AC}R_{BC})^3} \\
& \times f_{damp,zero}^{(9)}(R_{AB}, R_{AC}, R_{BC}).
\end{aligned} \tag{38}$$

The term in the second line is the three-body Axilrod–Teller–Muto (ATM) (What is this?????) term and the last line is the corresponding zero-damping function for this term.

$C_6^{AB}$  is the pairwise dipole-dipole dispersion coefficients calculated by numerical integration via the Casimir-Polder relation.

$$C_6^{AB} = \frac{3}{\pi} \sum_j w_j \bar{\alpha}_A(i\omega_j, q_A, CN_{cov}^A) \bar{\alpha}_B(i\omega_j, q_B, CN_{cov}^B) \tag{39}$$

$w_j$  are the integration weights, which are derived from a trapezoidal partitioning between the grid points  $j (j \in [1, 23])$ .

The isotropically averaged, dynamic dipole-dipole polarizabilities  $\bar{\alpha}$  at the  $j$ th imaginary frequency  $i\omega_j$  are obtained from the self-consistent D4 model; i.e., they are depending on the covalent coordination number and are also charge dependent.

$$\bar{\alpha}_A(i\omega_j, q_A, CN_{cov}^A) = \sum_r^{N_{A,ref}} \xi_A^r(q_A, q_{A,r}) \bar{\alpha}_{A,r}(i\omega_j, q_{A,r}, CN_{cov}^{A,r}) W_A^r(CN_{cov}^A, CN_{cov}^{A,r}) \tag{40}$$

The Gaussian weighting for each reference system is given by:

$$W_A^r(CN_{cov}^A, CN_{cov}^{A,r}) = \sum_{j=1}^{N_{gauss}} \frac{1}{\mathcal{N}} \exp[-6j \cdot (CN_{cov}^A - CN_{cov}^{A,r})^2] \tag{41}$$

with

$$\sum_r^{N_{A,ref}} W_A^r(CN_{cov}^A, CN_{cov}^{A,r}) = 1 \tag{42}$$

$\mathcal{N}$  is a normalization constant.

The number of Gaussian function per reference system  $N_{gauss}$  is mostly one, but equal to three for  $CN_{cov}^{A,r} = 0$  and reference systems with similar coordination number.

The charge-dependency is included via the empirical scaling function  $\xi_A^r$ .

$$\xi_A^r(q_A, q_{A,r}) = \exp \left[ 3 \left\{ 1 - \exp \left[ 4\eta_A \left( 1 - \frac{Z_A^{eff} + q_{A,r}}{Z_A^{eff} + q_A} \right) \right] \right\} \right] \quad (43)$$

where  $\eta_A$  is the chemical hardness taken from ref 98.

$Z_A^{eff}$  is the effective nuclear charge of atom A, which has been determined by subtracting the number of core electrons represented by the def2-ECPs in the time-dependent DFT reference calculations.

$C_8^{AB}$  is calculated recursively from the lowest order  $C_6^{AB}$  coefficients.

$$C_8^{AB} = 3C_6^{AB} \sqrt{Q^A Q^B} \quad (44)$$

$$Q^A = s_{42} \sqrt{Z^A} \frac{\langle r^4 \rangle^A}{\langle r^2 \rangle^A} \quad (45)$$

$\sqrt{Z^A}$  is the ad hoc nuclear charge dependent factor. (How do we get  $Z^{????}$ )

$\langle r^4 \rangle$  and  $\langle r^2 \rangle$  are simple multipole-type expectation values derived from atomic densities which are averaged geometrically to get the pair coefficients. (What is 'r', how we get $^{????}$ ) (what is  $s_{42}^{??}$ )

$C_{cov}^A$  is the covalent coordination number for atom A.

$q$  is the atomic charge, so  $q_A$  is the atomic charge for atom A.

The damping and scaling parameters in the dispersion model are:

$$a1 = 0.52 \quad | \quad a2 = 5.0 \quad | \quad s6 = 1.0 \quad | \quad s8 = 2.7 \quad | \quad s9 = 5.0$$

$C_9^{ABC}$  is the triple-dipole constant<sup>1</sup>:

$$C_9^{ABC} = \frac{3}{\pi} \int_0^\infty \alpha^A(i\omega) \alpha^B(i\omega) \alpha^C(i\omega) d\omega \quad (46)$$

The three-body contribution is typically  $< 5-10\%$  of  $E_{disp}$ , so it is small enough that we can reasonably approximate the coefficients by a geometric mean as<sup>1</sup>:

$$C_9^{ABC} \approx -\sqrt{C_6^{AB} C_6^{AC} C_6^{BC}} \quad (47)$$

$\theta_{ABC}$  is the angle between the two edges going from B to the other two atoms.

$\theta_{BCA}$  is the angle between the edges going from C to the other two and so on.

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<sup>1</sup>[https://www.researchgate.net/publication/43347348\\_A\\_Consistent\\_and\\_Accurate\\_Ab\\_Initio\\_Parametrization\\_of\\_Density\\_Functionals\\_for\\_the\\_94\\_Elements\\_H-Pu](https://www.researchgate.net/publication/43347348_A_Consistent_and_Accurate_Ab_Initio_Parametrization_of_Density_Functionals_for_the_94_Elements_H-Pu)

BJ = Becke-Johnson

$$f_n^{damp,BJ}(R_{AB}) = \frac{R_{AB}^n}{R_{AB}^n + (a_1 \cdot R_{AB}^{crit} + a_2)^6} \quad (48)$$

$$R_{AB}^{crit} = \sqrt{\frac{C_8^{AB}}{C_6^{AB}}} \quad (49)$$

$$f_9^{damp,zero}(R_{AB}, R_{AC}, R_{BC}) = \left( 1 + 6 \left( \sqrt{\frac{R_{AB}^{crit} R_{BC}^{crit} R_{CA}^{crit}}{R_{AB} R_{BC} R_{CA}}} \right)^{16} \right)^{-1} \quad (50)$$