



# wwPDB X-ray Structure Validation Summary Report

Mar 1, 2014 – 02:08 AM GMT

PDB ID : 1X7R  
Title : CRYSTAL STRUCTURE OF ESTROGEN RECEPTOR ALPHA COM-  
PLEXED WITH GENISTEIN  
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Deposited on : 2004-08-16  
Resolution : 2.00 Å(reported)

This is a wwPDB validation summary report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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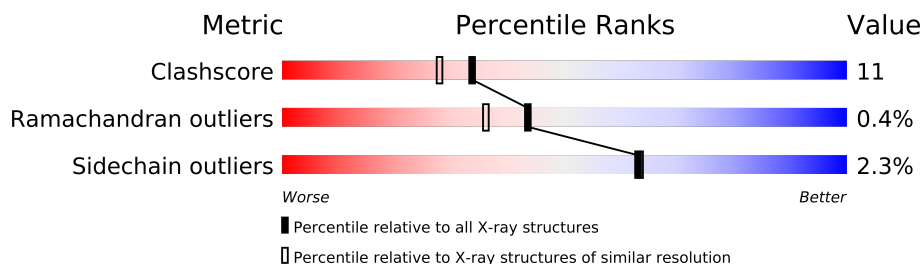
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 21963  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



| Metric                | Whole archive<br>(#Entries) | Similar resolution<br>(#Entries, resolution range(Å)) |
|-----------------------|-----------------------------|---|
| Clashscore            | 79885                       | 6188 (2.00-2.00)                                      |
| Ramachandran outliers | 78287                       | 6102 (2.00-2.00)                                      |
| Sidechain outliers    | 78261                       | 6100 (2.00-2.00)                                      |

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

| Mol | Chain | Length | Quality of chain |
|-----|-------|--------|------------------|
| 1   | A     | 245    |                  |
| 2   | B     | 13     |                  |

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

| Mol | Type | Chain | Res | Geometry | Electron density |
|-----|------|-------|-----|----------|------------------|
| 3   | GEN  | A     | 201 | X        | -                |

## 2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 2087 atoms, of which 0 are hydrogen and 0 are deuterium.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

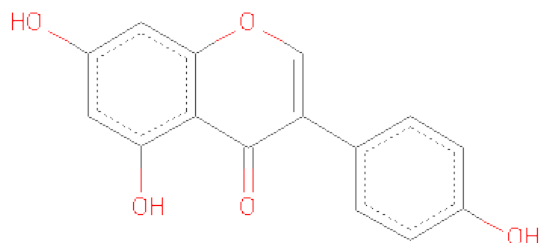
- Molecule 1 is a protein called Estrogen receptor 1 (alpha).

| Mol | Chain | Residues | Atoms |      |     |     |    | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|----|---------|---------|-------|
| 1   | A     | 235      | Total | C    | N   | O   | S  | 0       | 0       | 0     |
|     |       |          | 1871  | 1200 | 317 | 335 | 19 |         |         |       |

- Molecule 2 is a protein called steroid receptor coactivator-3.

| Mol | Chain | Residues | Atoms |    |    |    | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|----|----|----|---------|---------|-------|
| 2   | B     | 9        | Total | C  | N  | O  | 0       | 0       | 0     |
|     |       |          | 81    | 53 | 18 | 10 |         |         |       |

- Molecule 3 is GENISTEIN (three-letter code: GEN) (formula: C<sub>15</sub>H<sub>10</sub>O<sub>5</sub>).



| Mol | Chain | Residues | Atoms |    |   | ZeroOcc | AltConf |
|-----|-------|----------|-------|----|---|---------|---------|
| 3   | A     | 1        | Total | C  | O | 0       | 0       |
|     |       |          | 20    | 15 | 5 |         |         |

- Molecule 4 is water.

| Mol | Chain | Residues | Atoms        |          | ZeroOcc | AltConf |
|-----|-------|----------|--------------|----------|---------|---------|
| 4   | A     | 111      | Total<br>111 | O<br>111 | 0       | 0       |
| 4   | B     | 4        | Total<br>4   | O<br>4   | 0       | 0       |



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

| Property   | Value  | Source    |
|--|--|-----------|
| Space group  | P 21 21 2                                    | Depositor |
| Cell constants<br>a, b, c, $\alpha$ , $\beta$ , $\gamma$ | 60.88Å 69.83Å 57.37Å<br>90.00° 90.00° 90.00° | Depositor |
| Resolution (Å)   | 14.97 – 2.00                                 | Depositor |
| % Data completeness<br>(in resolution range)             | 92.6 (14.97-2.00)                            | Depositor |
| $R_{merge}$  | 0.05   | Depositor |
| $R_{sym}$  | 0.05   | Depositor |
| Refinement program                                       | CNS 1.1                                      | Depositor |
| R, $R_{free}$  | 0.224 , 0.269                                | Depositor |
| Estimated twinning fraction                              | No twinning to report.                       | Xtriage   |
| Total number of atoms                                    | 2087   | wwPDB-VP  |
| Average B, all atoms (Å <sup>2</sup> )                   | 28.0   | wwPDB-VP  |

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GEN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

| Mol | Chain | Bond lengths |             | Bond angles |             |
|-----|-------|--------------|-------------|-------------|-------------|
|     |       | RMSZ         | $\# Z  > 5$ | RMSZ        | $\# Z  > 5$ |
| 1   | A     | 0.32         | 0/1903      | 0.56        | 0/2570      |
| 2   | B     | 0.27         | 0/82        | 0.44        | 0/108       |
| All | All   | 0.32         | 0/1985      | 0.55        | 0/2678      |

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

| Mol | Chain | Non-H | H(model) | H(added) | Clashes | Symm-Clashes |
|-----|-------|-------|----------|----------|---------|--------------|
| 1   | A     | 1871  | 0        | 1932     | 42      | 0            |
| 2   | B     | 81    | 0        | 91       | 5       | 0            |
| 3   | A     | 20    | 0        | 10       | 1       | 0            |
| 4   | A     | 111   | 0        | 0        | 5       | 0            |
| 4   | B     | 4     | 0        | 0        | 1       | 0            |
| All | All   | 2087  | 0        | 2033     | 44      | 0            |

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 11.

The worst 5 of 44 close contacts within the same asymmetric unit are listed below.

| Atom-1           | Atom-2           | Distance(Å) | Clash(Å) |
|------------------|------------------|-------------|----------|
| 1:A:400:VAL:HG11 | 4:A:78:HOH:O     | 1.44        | 1.16     |
| 1:A:400:VAL:CG1  | 4:A:78:HOH:O     | 2.03        | 0.94     |
| 1:A:385:GLU:HG3  | 1:A:514:ILE:HG22 | 1.55        | 0.87     |
| 1:A:308:LEU:HA   | 1:A:481:LYS:HD2  | 1.63        | 0.81     |
| 1:A:465:THR:HG22 | 1:A:468:SER:H    | 1.46        | 0.79     |

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

| Mol | Chain | Analysed      | Favoured  | Allowed | Outliers | Percentiles |     |
|-----|-------|---------------|-----------|---------|----------|-------------|-----|
| 1   | A     | 231/245 (94%) | 224 (97%) | 6 (3%)  | 1 (0%)   | 43          | 36  |
| 2   | B     | 7/13 (54%)    | 7 (100%)  | 0       | 0        | 100         | 100 |
| All | All   | 238/258 (92%) | 231 (97%) | 6 (2%)  | 1 (0%)   | 43          | 36  |

All (1) Ramachandran outliers are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 1   | A     | 400 | VAL  |

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

| Mol | Chain | Analysed      | Rotameric | Outliers | Percentiles |    |
|-----|-------|---------------|-----------|----------|-------------|----|
| 1   | A     | 212/222 (96%) | 208 (98%) | 4 (2%)   | 69          | 71 |
| 2   | B     | 9/13 (69%)    | 8 (89%)   | 1 (11%)  | 9           | 5  |
| All | All   | 221/235 (94%) | 216 (98%) | 5 (2%)   | 63          | 63 |

All (5) residues with a non-rotameric sidechain are listed below:



| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 1   | A     | 365 | PRO  |
| 1   | A     | 396 | MET  |
| 1   | A     | 424 | ILE  |
| 1   | A     | 520 | LYS  |
| 2   | B     | 687 | HIS  |

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such sidechains are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 1   | A     | 474 | HIS  |
| 2   | B     | 695 | GLN  |
| 1   | A     | 519 | ASN  |
| 1   | A     | 441 | GLN  |
| 1   | A     | 502 | GLN  |

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

| Mol | Type | Chain | Res | Link | Bond lengths |      |          | Bond angles |      |          |
|-----|------|-------|-----|------|--------------|------|----------|-------------|------|----------|
|     |      |       |     |      | Counts       | RMSZ | # Z  > 2 | Counts      | RMSZ | # Z  > 2 |
| 3   | GEN  | A     | 201 | -    | 22,22,22     | 2.88 | 13 (59%) | 30,32,32    | 2.55 | 14 (46%) |

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

| Mol | Type | Chain | Res | Link | Chirals | Torsions | Rings   |
|-----|------|-------|-----|------|---------|----------|---------|
| 3   | GEN  | A     | 201 | -    | -       | 1/4/4/4  | 0/1/3/3 |

The worst 5 of 13 bond length outliers are listed below:

| Mol | Chain | Res | Type | Atoms   | Z    | Observed(Å) | Ideal(Å) |
|-----|-------|-----|------|---------|------|-------------|----------|
| 3   | A     | 201 | GEN  | C6-C5   | 6.24 | 1.50        | 1.40     |
| 3   | A     | 201 | GEN  | C6-C7   | 4.52 | 1.53        | 1.39     |
| 3   | A     | 201 | GEN  | O9-C8   | 4.03 | 1.42        | 1.34     |
| 3   | A     | 201 | GEN  | C7-C11  | 3.95 | 1.56        | 1.50     |
| 3   | A     | 201 | GEN  | C16-C15 | 3.84 | 1.46        | 1.38     |

The worst 5 of 14 bond angle outliers are listed below:

| Mol | Chain | Res | Type | Atoms       | Z     | Observed(°) | Ideal(°) |
|-----|-------|-----|------|-------------|-------|-------------|----------|
| 3   | A     | 201 | GEN  | O9-C10-C5   | 4.44  | 125.84      | 121.29   |
| 3   | A     | 201 | GEN  | C7-C6-C5    | -4.40 | 114.30      | 121.04   |
| 3   | A     | 201 | GEN  | C16-C15-C14 | -4.13 | 114.89      | 119.87   |
| 3   | A     | 201 | GEN  | C1-C2-C3    | 3.98  | 124.38      | 120.91   |
| 3   | A     | 201 | GEN  | C16-C11-C12 | 3.86  | 124.78      | 117.58   |

There are no chirality outliers.

All (1) torsion outliers are listed below:

| Mol | Chain | Res | Type | Atoms         |
|-----|-------|-----|------|---------------|
| 3   | A     | 201 | GEN  | C12-C11-C7-C6 |

There are no ring outliers.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.