



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 13, 2026 – 06:46 PM UTC

PDB ID : 3EK7 / pdb\_00003ek7  
Title : Calcium-saturated GCaMP2 dimer  
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Deposited on : 2008-09-18  
Resolution : 1.85 Å(reported)

This is a wwPDB X-ray Structure 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

|                                |   |                                                                  |
|--------------------------------|---|------------------------------------------------------------------|
| MolProbity                     | : | 4-5-2 with Phenix2.0                                             |
| Mogul                          | : | 2022.3.0, CSD as543be (2022)                                     |
| Xtriage (Phenix)               | : | 2.0                                                              |
| EDS                            | : | <b>FAILED</b>                                                    |
| Percentile statistics          | : | 20250101.v01 (using entries in the PDB archive January 1st 2025) |
| Ideal geometry (proteins)      | : | Engh & Huber (2001)                                              |
| Ideal geometry (DNA, RNA)      | : | Parkinson et al. (1996)                                          |
| Validation Pipeline (wwPDB-VP) | : | 2.49                                                             |

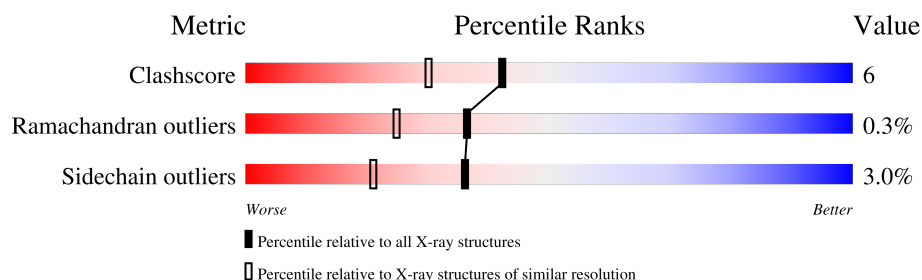
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.85 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            | 190562                      | 3579 (1.86-1.86)                                      |
| Ramachandran outliers | 187476                      | 3553 (1.86-1.86)                                      |
| Sidechain outliers    | 187428                      | 3553 (1.86-1.86)                                      |

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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ .

Note EDS failed to run properly.

| Mol | Chain | Length | Quality of chain  |
|-----|-------|--------|-------------------|
| 1   | A     | 449    | <br>75% 10% • 13% |

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3333 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Myosin light chain kinase, Green fluorescent protein, Calmodulin chimera.

| Mol | Chain | Residues | Atoms |      |     |     |    | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|----|---------|---------|-------|
| 1   | A     | 390      | Total | C    | N   | O   | S  | 0       | 4       | 0     |
|     |       |          | 3131  | 1966 | 526 | 623 | 16 |         |         |       |

There are 20 discrepancies between the modelled and reference sequences:

| Chain | Residue | Modelled | Actual | Comment        | Reference  |
|-------|---------|----------|--------|----------------|------------|
| A     | 76      | ALA      | VAL    | see remark 999 | UNP P42212 |
| A     | 88      | GLY      | SER    | see remark 999 | UNP P42212 |
| A     | 93      | TYR      | ASP    | see remark 999 | UNP P42212 |
| A     | 119     | LYS      | ALA    | see remark 999 | UNP P42212 |
| A     | 144     | LEU      | HIS    | see remark 999 | UNP P42212 |
| A     | 152     | GLY      | -      | linker         | UNP P42212 |
| A     | 153     | GLY      | -      | linker         | UNP P42212 |
| A     | 154     | THR      | -      | linker         | UNP P42212 |
| A     | 155     | GLY      | -      | linker         | UNP P42212 |
| A     | 156     | GLY      | -      | linker         | UNP P42212 |
| A     | 157     | SER      | -      | linker         | UNP P42212 |
| A     | 158     | MET      | -      | linker         | UNP P42212 |
| A     | 159     | VAL      | -      | linker         | UNP P42212 |
| A     | 222     | LEU      | PHE    | see remark 999 | UNP P42212 |
| A     | 224     | CRO      | SER    | chromophore    | UNP P42212 |
| A     | ?       | -        | TYR    | chromophore    | UNP P42212 |
| A     | ?       | -        | GLY    | chromophore    | UNP P42212 |
| A     | 251     | ILE      | VAL    | see remark 999 | UNP P42212 |
| A     | 303     | THR      | -      | linker         | UNP P42212 |
| A     | 304     | ARG      | -      | linker         | UNP P42212 |

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

| Mol | Chain | Residues | Atoms |    | ZeroOcc | AltConf |
|-----|-------|----------|-------|----|---------|---------|
| 2   | A     | 4        | Total | Ca | 0       | 0       |
|     |       |          | 4     | 4  |         |         |

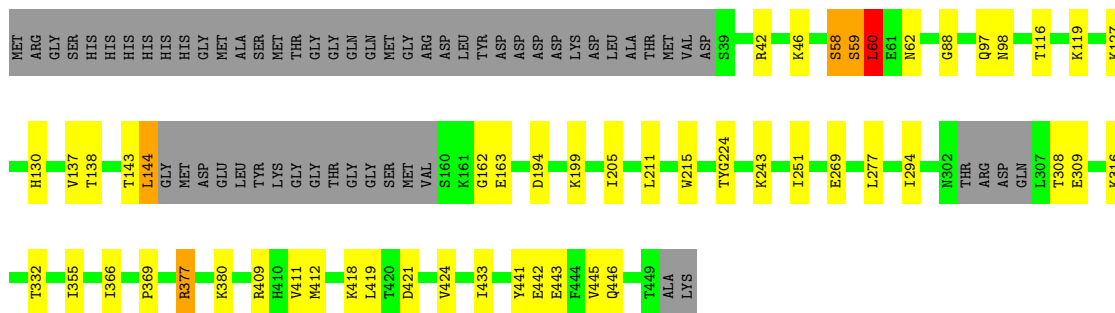
- Molecule 3 is water.

| Mol | Chain | Residues | Atoms        |          | ZeroOcc | AltConf |
|-----|-------|----------|--------------|----------|---------|---------|
| 3   | A     | 198      | Total<br>198 | O<br>198 | 0       | 0       |



Note EDS failed to run properly.

- Molecule 1: Myosin light chain kinase, Green fluorescent protein, Calmodulin chimera



## 4 Data and refinement statistics

EDS failed to run properly - this section is therefore incomplete.

| Property                                                 | Value                                                       | Source    |
|----------------------------------------------------------|-------------------------------------------------------------|-----------|
| Space group                                              | C 1 2 1                                                     | Depositor |
| Cell constants<br>a, b, c, $\alpha$ , $\beta$ , $\gamma$ | 126.13Å 47.31Å 68.94Å<br>90.00° 100.48° 90.00°              | Depositor |
| Resolution (Å)                                           | 20.67 – 1.85                                                | Depositor |
| % Data completeness<br>(in resolution range)             | 98.5 (20.67-1.85)                                           | Depositor |
| $R_{merge}$                                              | (Not available)                                             | Depositor |
| $R_{sym}$                                                | 0.08                                                        | Depositor |
| $\langle I/\sigma(I) \rangle$ <sup>1</sup>               | 5.26 (at 1.85Å)                                             | Xtriage   |
| Refinement program                                       | REFMAC 5.2.0019                                             | Depositor |
| R, $R_{free}$                                            | 0.189 , 0.241                                               | Depositor |
| Wilson B-factor (Å <sup>2</sup> )                        | 21.6                                                        | Xtriage   |
| Anisotropy                                               | 0.329                                                       | Xtriage   |
| L-test for twinning <sup>2</sup>                         | $\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$ | Xtriage   |
| Estimated twinning fraction                              | No twinning to report.                                      | Xtriage   |
| Total number of atoms                                    | 3333                                                        | wwPDB-VP  |
| Average B, all atoms (Å <sup>2</sup> )                   | 25.0                                                        | wwPDB-VP  |

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.56% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

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

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     | 1.05         | 4/3165 (0.1%) | 0.99        | 3/4259 (0.1%) |

All (4) bond length outliers are listed below:

| Mol | Chain | Res | Type | Atoms  | Z    | Observed(Å) | Ideal(Å) |
|-----|-------|-----|------|--------|------|-------------|----------|
| 1   | A     | 366 | ILE  | CA-CB  | 6.58 | 1.61        | 1.54     |
| 1   | A     | 62  | ASN  | CG-OD1 | 6.07 | 1.35        | 1.23     |
| 1   | A     | 446 | GLN  | CD-OE1 | 5.69 | 1.34        | 1.23     |
| 1   | A     | 97  | GLN  | CD-OE1 | 5.08 | 1.33        | 1.23     |

All (3) bond angle outliers are listed below:

| Mol | Chain | Res | Type | Atoms  | Z     | Observed(°) | Ideal(°) |
|-----|-------|-----|------|--------|-------|-------------|----------|
| 1   | A     | 60  | LEU  | N-CA-C | -7.38 | 96.02       | 108.26   |
| 1   | A     | 88  | GLY  | N-CA-C | -6.05 | 105.87      | 112.04   |
| 1   | A     | 58  | SER  | N-CA-C | 5.05  | 117.68      | 111.82   |

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-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 hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

| Mol | Chain | Non-H | H(model) | H(added) | Clashes | Symm-Clashes |
|-----|-------|-------|----------|----------|---------|--------------|
| 1   | A     | 3131  | 0        | 3034     | 39      | 0            |

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| Mol | Chain | Non-H | H(model) | H(added) | Clashes | Symm-Clashes |
|-----|-------|-------|----------|----------|---------|--------------|
| 2   | A     | 4     | 0        | 0        | 0       | 0            |
| 3   | A     | 198   | 0        | 0        | 4       | 1            |
| All | All   | 3333  | 0        | 3034     | 39      | 1            |

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

The worst 5 of 39 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

| Atom-1           | Atom-2           | Interatomic distance (Å) | Clash overlap (Å) |
|------------------|------------------|--------------------------|-------------------|
| 1:A:409:ARG:HG3  | 1:A:424:VAL:HG21 | 1.51                     | 0.89              |
| 1:A:194:ASP:OD2  | 3:A:611:HOH:O    | 1.95                     | 0.84              |
| 1:A:143:THR:HG23 | 1:A:144:LEU:HD12 | 1.62                     | 0.81              |
| 1:A:163:GLU:HG3  | 3:A:527:HOH:O    | 1.88                     | 0.73              |
| 1:A:412:MET:HE1  | 1:A:419:LEU:HD11 | 1.71                     | 0.71              |

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

| Atom-1        | Atom-2               | Interatomic distance (Å) | Clash overlap (Å) |
|---------------|----------------------|--------------------------|-------------------|
| 3:A:627:HOH:O | 3:A:643:HOH:O[1_545] | 1.76                     | 0.44              |

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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     | 385/449 (86%) | 380 (99%) | 4 (1%)  | 1 (0%)   | 36 25       |

All (1) Ramachandran outliers are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 1   | A     | 377 | ARG  |

### 5.3.2 Protein sidechains [i](#)

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     | 340/382 (89%) | 329 (97%) | 11 (3%)  | 34 19       |

5 of 11 residues with a non-rotameric sidechain are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 1   | A     | 144 | LEU  |
| 1   | A     | 308 | THR  |
| 1   | A     | 380 | LYS  |
| 1   | A     | 309 | GLU  |
| 1   | A     | 60  | LEU  |

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 1   | A     | 77  | ASN  |
| 1   | A     | 279 | ASN  |
| 1   | A     | 293 | ASN  |
| 1   | A     | 345 | ASN  |
| 1   | A     | 438 | GLN  |

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

1 non-standard protein/DNA/RNA residue 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$ |
| 1   | CRO  | A     | 224 | 1    | 22,23,24     | 3.28 | 5 (22%)     | 30,32,34    | 2.61 | 8 (26%)     |

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   |
|-----|------|-------|-----|------|---------|------------|---------|
| 1   | CRO  | A     | 224 | 1    | -       | 1/12/31/32 | 0/2/2/2 |

All (5) bond length outliers are listed below:

| Mol | Chain | Res | Type | Atoms   | Z     | Observed(Å) | Ideal(Å) |
|-----|-------|-----|------|---------|-------|-------------|----------|
| 1   | A     | 224 | CRO  | CB2-CA2 | 12.65 | 1.47        | 1.35     |
| 1   | A     | 224 | CRO  | CA2-C2  | -5.44 | 1.42        | 1.48     |
| 1   | A     | 224 | CRO  | C1-N2   | 4.01  | 1.38        | 1.32     |
| 1   | A     | 224 | CRO  | C2-N3   | -3.76 | 1.31        | 1.40     |
| 1   | A     | 224 | CRO  | O2-C2   | 2.83  | 1.28        | 1.23     |

The worst 5 of 8 bond angle outliers are listed below:

| Mol | Chain | Res | Type | Atoms     | Z     | Observed(°) | Ideal(°) |
|-----|-------|-----|------|-----------|-------|-------------|----------|
| 1   | A     | 224 | CRO  | O2-C2-CA2 | -9.33 | 125.06      | 131.02   |
| 1   | A     | 224 | CRO  | CA2-C2-N3 | 6.39  | 108.87      | 103.50   |
| 1   | A     | 224 | CRO  | C2-N3-C1  | -3.62 | 106.39      | 108.07   |
| 1   | A     | 224 | CRO  | C1-CA1-N1 | -3.18 | 104.09      | 109.78   |
| 1   | A     | 224 | CRO  | CA2-N2-C1 | -2.92 | 103.52      | 105.80   |

There are no chirality outliers.

All (1) torsion outliers are listed below:

| Mol | Chain | Res | Type | Atoms         |
|-----|-------|-----|------|---------------|
| 1   | A     | 224 | CRO  | N3-C1-CA1-CB1 |

There are no ring outliers.

No monomer is involved in short contacts.

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

EDS failed to run properly - this section is therefore empty.

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

EDS failed to run properly - this section is therefore empty.

### 6.3 Carbohydrates

EDS failed to run properly - this section is therefore empty.

### 6.4 Ligands

EDS failed to run properly - this section is therefore empty.

### 6.5 Other polymers

EDS failed to run properly - this section is therefore empty.