



# Full wwPDB NMR Structure Validation Report ⓘ

Mar 6, 2026 – 10:41 PM UTC

PDB ID : 2G5M / pdb\_00002g5m  
Title : Spinophilin PDZ domain  
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Deposited on : 2006-02-23

This is a Full wwPDB NMR Structure Validation 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/NMRValidationReportHelp>

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                                             |
| Percentile statistics          | : | 20250101.v01 (using entries in the PDB archive January 1st 2025) |
| wwPDB-RCI                      | : | v_1n_11_5_13_A (Berjanski et al., 2005)                          |
| PANAV                          | : | Wang et al. (2010)                                               |
| wwPDB-ShiftChecker             | : | v1.2                                                             |
| Ideal geometry (proteins)      | : | Engh & Huber (2001)                                              |
| Ideal geometry (DNA, RNA)      | : | Parkinson et al. (1996)                                          |
| Validation Pipeline (wwPDB-VP) | : | 2.49                                                             |

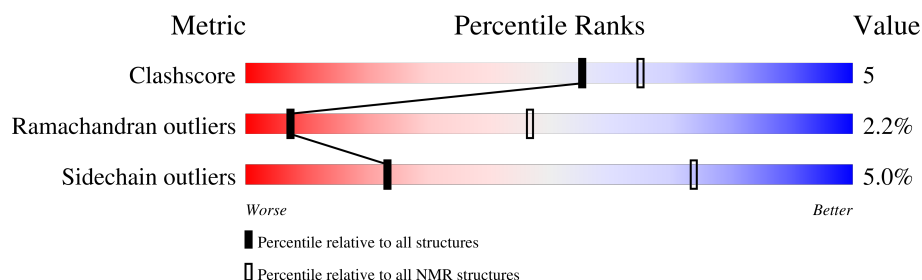
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*SOLUTION NMR*

The overall completeness of chemical shifts assignment was not calculated.

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) | NMR archive<br>(#Entries) |
|-----------------------|-----------------------------|---------------------------|
| Clashscore            | 229148                      | 14424                     |
| Ramachandran outliers | 224038                      | 12848                     |
| Sidechain outliers    | 223484                      | 12823                     |

The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A cyan segment indicates the fraction of residues that are not part of the well-defined cores, and 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  $\leq 5\%$

| Mol | Chain | Length | Quality of chain |
|-----|-------|--------|------------------|
| 1   | B     | 113    |                  |

## 2 Ensemble composition and analysis

This entry contains 20 models. Model 11 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *closest to the average*.

The following residues are included in the computation of the global validation metrics.

| Well-defined (core) protein residues |                          |                   |              |
|--------------------------------------|--------------------------|-------------------|--------------|
| Well-defined core                    | Residue range (total)    | Backbone RMSD (Å) | Medoid model |
| 1                                    | B:3-B:24, B:35-B:93 (81) | 0.74              | 11           |

Ill-defined regions of proteins are excluded from the global statistics.

Ligands and non-protein polymers are included in the analysis.

The models can be grouped into 5 clusters and 2 single-model clusters were found.

| Cluster number        | Models                     |
|-----------------------|----------------------------|
| 1                     | 2, 6, 8, 9, 11, 15, 18, 19 |
| 2                     | 5, 7, 12                   |
| 3                     | 1, 13, 14                  |
| 4                     | 10, 17                     |
| 5                     | 3, 4                       |
| Single-model clusters | 16; 20                     |

### 3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 1706 atoms, of which 857 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Neurabin-2.

| Mol | Chain | Residues | Atoms |     |     |     |     |   | Trace |
|-----|-------|----------|-------|-----|-----|-----|-----|---|-------|
| 1   | B     | 113      | Total | C   | H   | N   | O   | S | 0     |
|     |       |          | 1706  | 525 | 857 | 151 | 169 | 4 |       |

There are 3 discrepancies between the modelled and reference sequences:

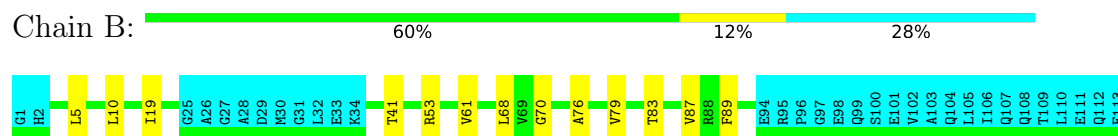
| Chain | Residue | Modelled | Actual | Comment          | Reference  |
|-------|---------|----------|--------|------------------|------------|
| B     | 1       | GLY      | -      | cloning artifact | UNP O35274 |
| B     | 2       | HIS      | -      | cloning artifact | UNP O35274 |
| B     | 3       | MET      | -      | cloning artifact | UNP O35274 |

## 4 Residue-property plots

### 4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outliers are shown as green connectors. Residues which are classified as ill-defined in the NMR ensemble, are shown in cyan with an underline colour-coded according to the previous scheme. Residues which were present in the experimental sample, but not modelled in the final structure are shown in grey.

- Molecule 1: Neurabin-2

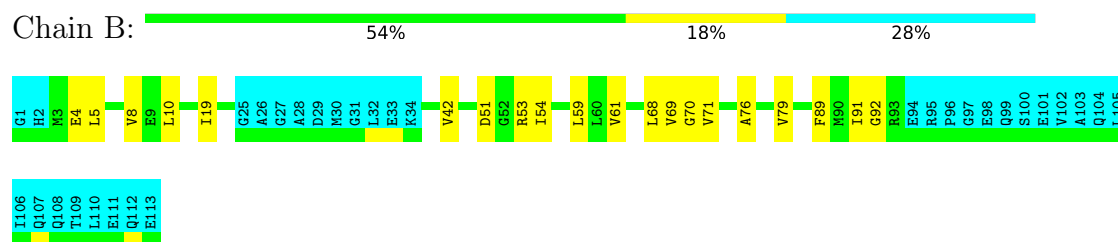


### 4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

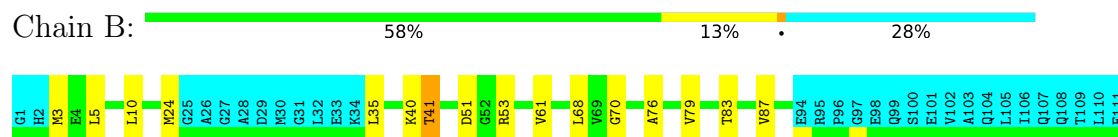
#### 4.2.1 Score per residue for model 1

- Molecule 1: Neurabin-2



#### 4.2.2 Score per residue for model 2

- Molecule 1: Neurabin-2



Q112  
E113

### 4.2.3 Score per residue for model 3

- Molecule 1: Neurabin-2

Chain B:  56% 14% • 28%

G1 H2 P7 L10 L19 G25 A26 G27 A28 D29 M30 G31 L32 E33 K34 E44 A48 H49 V61 E62 V63 A76 V79 T83 K84 G85 R86 V87 R88 F89 M90 I91 G92 R93 E94 R95 P96 G97 G98 E99 Q99 S100 V102 V103 A103 Q104 Q105 L106 I107 Q108 T109

L110 E111 Q112 E113

### 4.2.4 Score per residue for model 4

- Molecule 1: Neurabin-2

Chain B:  58% 12% • 28%

G1 H2 L5 L10 L19 G25 A26 G27 A28 D29 M30 G31 L32 E33 K34 T41 A48 H49 R50 V61 L68 V69 G70 A76 V79 T83 R86 V87 R93 E94 R95 P96 G97 E98 Q99 S100 V102 V103 A103 Q104 L105 I106 Q107 Q108 T109 E111

Q112  
E113

### 4.2.5 Score per residue for model 5

- Molecule 1: Neurabin-2

Chain B:  56% 15% • 28%

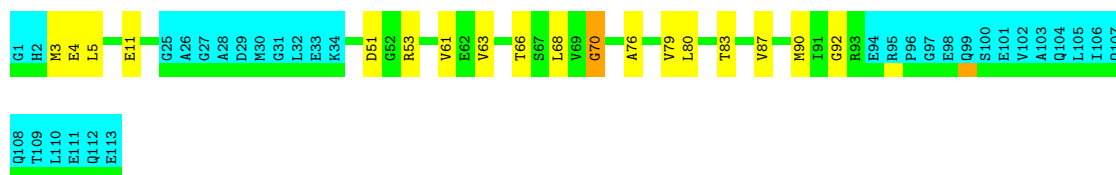
G1 H2 L5 V8 E9 L10 D13 G25 A26 G27 A28 D29 M30 G31 L32 E33 R34 L35 A47 R53 I54 V61 L68 V69 G70 A76 V79 T83 F89 M90 I91 G92 R93 E94 R95 P96 G97 G98 E99 Q99 S100 V102 V103 A103 Q104 Q105 L106 I107 Q108

T109 L110 E111 Q112 E113

### 4.2.6 Score per residue for model 6

- Molecule 1: Neurabin-2

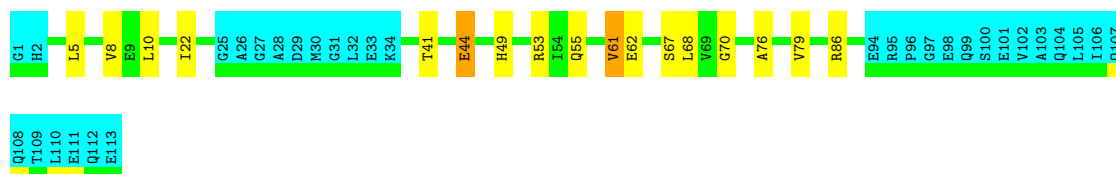
Chain B:  56% 15% • 28%



#### 4.2.7 Score per residue for model 7

- Molecule 1: Neurabin-2

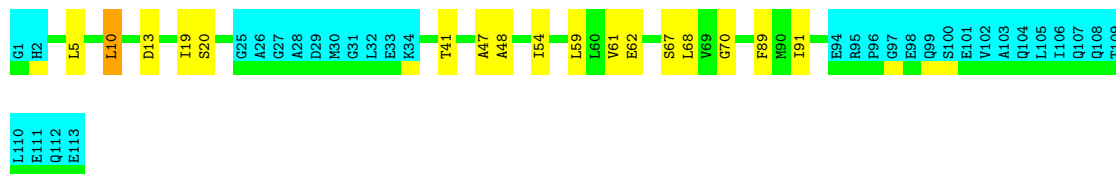
Chain B: 57% 13% 28%



#### 4.2.8 Score per residue for model 8

- Molecule 1: Neurabin-2

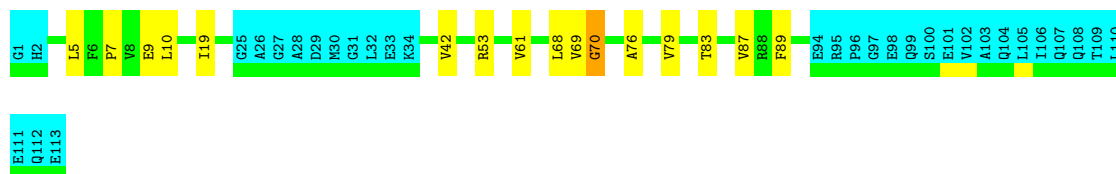
Chain B: 57% 14% 28%



#### 4.2.9 Score per residue for model 9

- Molecule 1: Neurabin-2

Chain B: 58% 13% 28%



#### 4.2.10 Score per residue for model 10

- Molecule 1: Neurabin-2

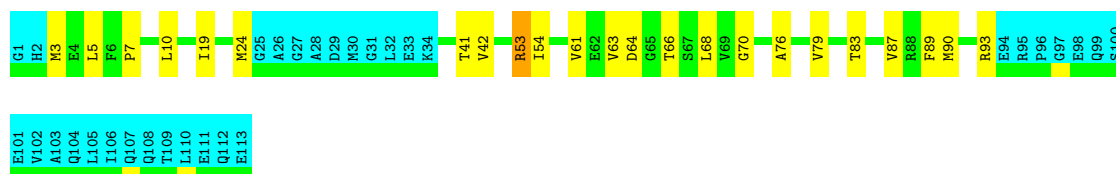
Chain B:  58% 11% • 28%



#### 4.2.11 Score per residue for model 11 (medoid)

- Molecule 1: Neurabin-2

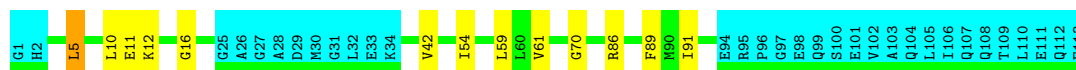
Chain B:  51% 19% • 28%



#### 4.2.12 Score per residue for model 12

- Molecule 1: Neurabin-2

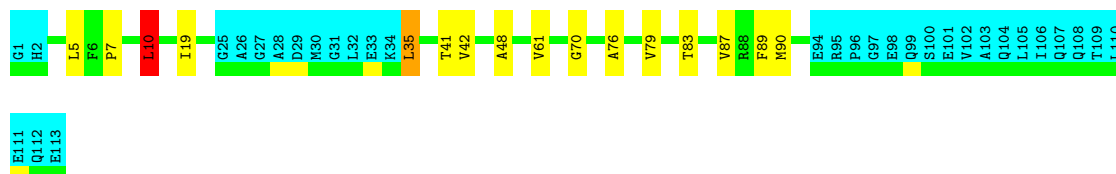
Chain B:  60% 11% • 28%



#### 4.2.13 Score per residue for model 13

- Molecule 1: Neurabin-2

Chain B:  58% 12% •• 28%



#### 4.2.14 Score per residue for model 14

- Molecule 1: Neurabin-2

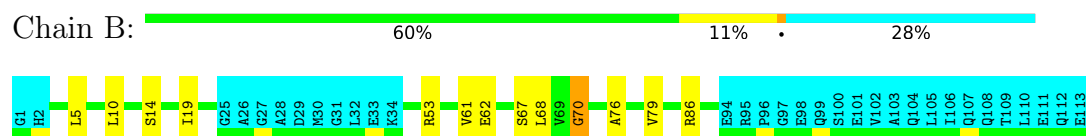
Chain B:  59% 12% 28%





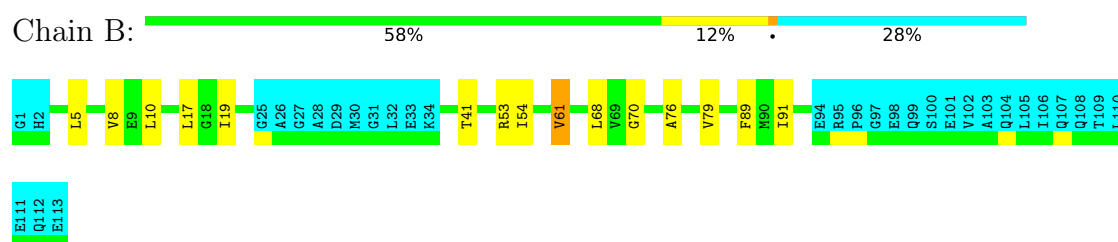
#### 4.2.15 Score per residue for model 15

- Molecule 1: Neurabin-2



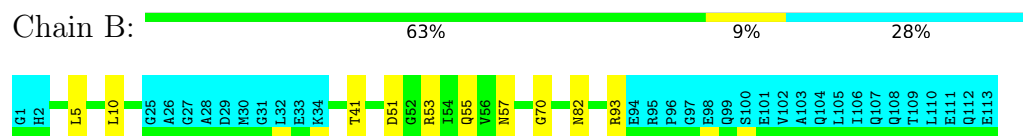
#### 4.2.16 Score per residue for model 16

- Molecule 1: Neurabin-2



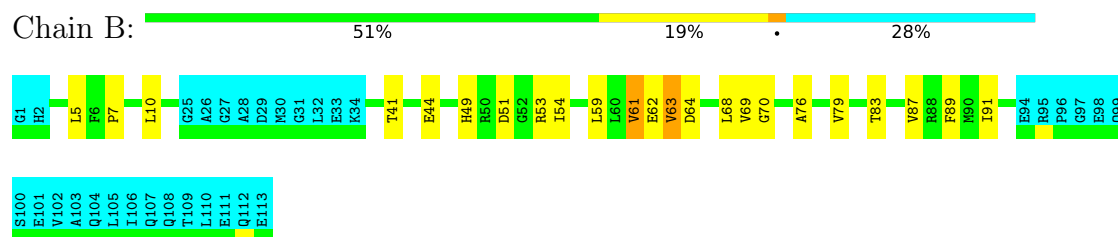
#### 4.2.17 Score per residue for model 17

- Molecule 1: Neurabin-2



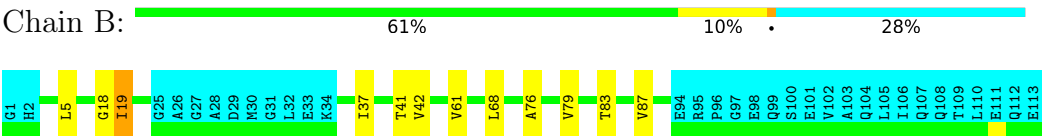
#### 4.2.18 Score per residue for model 18

- Molecule 1: Neurabin-2



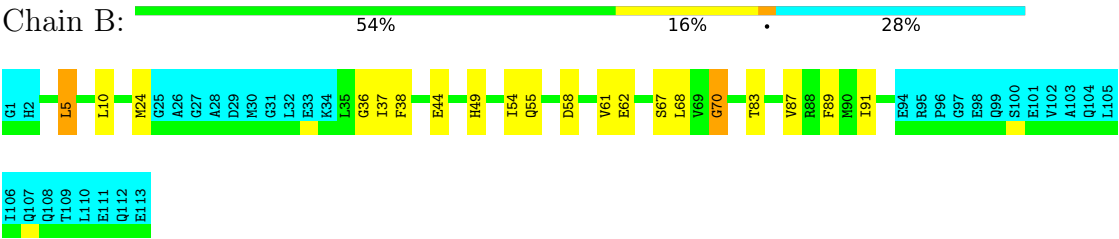
#### 4.2.19 Score per residue for model 19

- Molecule 1: Neurabin-2



4.2.20 Score per residue for model 20

- Molecule 1: Neurabin-2



## 5 Refinement protocol and experimental data overview

The models were refined using the following method: *simulated annealing torsion angle dynamics*.

Of the 50 calculated structures, 20 were deposited, based on the following criterion: *structures with the lowest energy*.

The following table shows the software used for structure solution, optimisation and refinement.

| Software name | Classification     | Version |
|---------------|--------------------|---------|
| CYANA         | structure solution | 2.0     |
| CNS           | refinement         | 1.1     |
| atnos/candid  | structure solution | 2.0     |

No chemical shift data was provided.

## 6 Model quality [i](#)

### 6.1 Standard geometry [i](#)

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 (average) 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   | B     | 1.27±0.06    | 1±1/615 ( 0.2± 0.1%) | 0.96±0.03   | 0±0/827 ( 0.0± 0.0%) |
| All | All   | 1.27         | 19/12300 ( 0.2%)     | 0.96        | 1/16540 ( 0.0%)      |

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

| Mol | Chain | Chirality | Planarity |
|-----|-------|-----------|-----------|
| 1   | B     | 0.0±0.0   | 0.3±0.7   |
| All | All   | 0         | 6         |

All unique bond outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

| Mol | Chain | Res | Type | Atoms | Z      | Observed(Å) | Ideal(Å) | Models |       |
|-----|-------|-----|------|-------|--------|-------------|----------|--------|-------|
|     |       |     |      |       |        |             |          | Worst  | Total |
| 1   | B     | 35  | LEU  | C-N   | -12.91 | 1.27        | 1.33     | 13     | 1     |
| 1   | B     | 41  | THR  | N-CA  | -6.83  | 1.37        | 1.45     | 2      | 1     |
| 1   | B     | 19  | ILE  | C-N   | -6.44  | 1.25        | 1.33     | 9      | 8     |
| 1   | B     | 22  | ILE  | C-N   | -5.54  | 1.30        | 1.33     | 7      | 1     |
| 1   | B     | 20  | SER  | C-N   | -5.41  | 1.27        | 1.33     | 8      | 1     |
| 1   | B     | 92  | GLY  | C-N   | -5.38  | 1.26        | 1.33     | 5      | 3     |
| 1   | B     | 61  | VAL  | C-N   | -5.26  | 1.27        | 1.33     | 12     | 1     |
| 1   | B     | 10  | LEU  | C-N   | -5.19  | 1.26        | 1.33     | 13     | 1     |
| 1   | B     | 61  | VAL  | C-O   | -5.15  | 1.18        | 1.24     | 16     | 1     |
| 1   | B     | 70  | GLY  | C-N   | -5.07  | 1.27        | 1.33     | 6      | 1     |

All unique angle outliers are listed below.

| Mol | Chain | Res | Type | Atoms    | Z    | Observed(°) | Ideal(°) | Models |       |
|-----|-------|-----|------|----------|------|-------------|----------|--------|-------|
|     |       |     |      |          |      |             |          | Worst  | Total |
| 1   | B     | 57  | ASN  | CA-CB-CG | 5.02 | 117.62      | 112.60   | 10     | 1     |

There are no chirality outliers.

All unique planar outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

| Mol | Chain | Res | Type | Group     | Models (Total) |
|-----|-------|-----|------|-----------|----------------|
| 1   | B     | 86  | ARG  | Sidechain | 3              |
| 1   | B     | 93  | ARG  | Sidechain | 2              |
| 1   | B     | 50  | ARG  | Sidechain | 1              |

## 6.2 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 each 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 averaged over the ensemble.

| Mol | Chain | Non-H | H(model) | H(added) | Clashes |
|-----|-------|-------|----------|----------|---------|
| 1   | B     | 609   | 631      | 631      | 6±2     |
| All | All   | 12180 | 12620    | 12620    | 124     |

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 5.

All unique clashes are listed below, sorted by their clash magnitude.

| Atom-1          | Atom-2          | Clash(Å) | Distance(Å) | Models |       |
|-----------------|-----------------|----------|-------------|--------|-------|
|                 |                 |          |             | Worst  | Total |
| 1:B:55:GLN:HB2  | 1:B:58:ASP:OD1  | 0.62     | 1.93        | 20     | 1     |
| 1:B:61:VAL:O    | 1:B:68:LEU:HB2  | 0.62     | 1.94        | 7      | 15    |
| 1:B:83:THR:HB   | 1:B:87:VAL:HG11 | 0.59     | 1.74        | 10     | 12    |
| 1:B:76:ALA:O    | 1:B:79:VAL:HG22 | 0.58     | 1.98        | 14     | 2     |
| 1:B:76:ALA:O    | 1:B:79:VAL:HB   | 0.56     | 2.01        | 13     | 1     |
| 1:B:24:MET:HB2  | 1:B:38:PHE:CD1  | 0.55     | 2.35        | 20     | 1     |
| 1:B:54:ILE:HD11 | 1:B:89:PHE:CE2  | 0.53     | 2.38        | 16     | 9     |
| 1:B:53:ARG:HA   | 1:B:53:ARG:NE   | 0.53     | 2.19        | 10     | 3     |
| 1:B:51:ASP:OD2  | 1:B:53:ARG:HD3  | 0.53     | 2.03        | 6      | 3     |
| 1:B:8:VAL:HG13  | 1:B:53:ARG:HB3  | 0.51     | 1.81        | 7      | 3     |
| 1:B:59:LEU:O    | 1:B:91:ILE:HA   | 0.51     | 2.05        | 18     | 5     |
| 1:B:61:VAL:HB   | 1:B:90:MET:O    | 0.50     | 2.06        | 6      | 6     |
| 1:B:8:VAL:HG13  | 1:B:53:ARG:HG3  | 0.49     | 1.84        | 1      | 1     |
| 1:B:51:ASP:OD1  | 1:B:53:ARG:HG2  | 0.49     | 2.07        | 10     | 3     |
| 1:B:76:ALA:O    | 1:B:80:LEU:HG   | 0.49     | 2.07        | 10     | 2     |
| 1:B:7:PRO:HA    | 1:B:89:PHE:O    | 0.49     | 2.07        | 13     | 5     |
| 1:B:35:LEU:O    | 1:B:70:GLY:HA2  | 0.49     | 2.08        | 13     | 2     |

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| Atom-1          | Atom-2          | Clash(Å) | Distance(Å) | Models |       |
|-----------------|-----------------|----------|-------------|--------|-------|
|                 |                 |          |             | Worst  | Total |
| 1:B:11:GLU:OE2  | 1:B:86:ARG:HD3  | 0.49     | 2.07        | 12     | 1     |
| 1:B:76:ALA:O    | 1:B:79:VAL:HG12 | 0.48     | 2.08        | 7      | 12    |
| 1:B:44:GLU:H    | 1:B:44:GLU:CD   | 0.48     | 2.17        | 7      | 1     |
| 1:B:44:GLU:HA   | 1:B:49:HIS:CG   | 0.47     | 2.44        | 20     | 3     |
| 1:B:4:GLU:O     | 1:B:92:GLY:HA2  | 0.46     | 2.10        | 6      | 3     |
| 1:B:63:VAL:HG12 | 1:B:66:THR:HB   | 0.45     | 1.87        | 6      | 2     |
| 1:B:13:ASP:HB2  | 1:B:47:ALA:N    | 0.45     | 2.26        | 5      | 1     |
| 1:B:62:GLU:CB   | 1:B:67:SER:HA   | 0.45     | 2.42        | 20     | 3     |
| 1:B:62:GLU:HG2  | 1:B:63:VAL:O    | 0.44     | 2.11        | 18     | 2     |
| 1:B:10:LEU:HD11 | 1:B:48:ALA:HB2  | 0.44     | 1.88        | 13     | 3     |
| 1:B:9:GLU:O     | 1:B:53:ARG:HG3  | 0.43     | 2.13        | 9      | 1     |
| 1:B:5:LEU:HA    | 1:B:91:ILE:O    | 0.42     | 2.14        | 20     | 2     |
| 1:B:24:MET:O    | 1:B:35:LEU:HG   | 0.42     | 2.15        | 2      | 1     |
| 1:B:64:ASP:HB2  | 1:B:87:VAL:HB   | 0.42     | 1.90        | 11     | 2     |
| 1:B:44:GLU:HA   | 1:B:49:HIS:HB3  | 0.42     | 1.92        | 3      | 1     |
| 1:B:13:ASP:HB2  | 1:B:47:ALA:HB2  | 0.42     | 1.91        | 8      | 1     |
| 1:B:68:LEU:HD12 | 1:B:71:VAL:CG2  | 0.41     | 2.45        | 1      | 1     |
| 1:B:79:VAL:O    | 1:B:83:THR:HG23 | 0.41     | 2.15        | 5      | 2     |
| 1:B:68:LEU:C    | 1:B:70:GLY:H    | 0.41     | 2.23        | 9      | 1     |
| 1:B:62:GLU:HB2  | 1:B:67:SER:HA   | 0.41     | 1.91        | 7      | 1     |
| 1:B:10:LEU:O    | 1:B:86:ARG:HG3  | 0.41     | 2.15        | 3      | 1     |
| 1:B:3:MET:HA    | 1:B:93:ARG:O    | 0.40     | 2.16        | 11     | 1     |
| 1:B:8:VAL:HB    | 1:B:91:ILE:HD13 | 0.40     | 1.94        | 16     | 1     |
| 1:B:4:GLU:OE2   | 1:B:93:ARG:HD2  | 0.40     | 2.16        | 10     | 1     |
| 1:B:10:LEU:HD21 | 1:B:48:ALA:HA   | 0.40     | 1.92        | 3      | 1     |
| 1:B:44:GLU:HA   | 1:B:49:HIS:CB   | 0.40     | 2.46        | 3      | 1     |

## 6.3 Torsion angles ⓘ

### 6.3.1 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 PDB entries followed by that with respect to all NMR entries. 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   | B     | 81/113 (72%)    | 69±3 (85±4%) | 10±3 (12±3%) | 2±1 (2±1%) | 7           | 47 |
| All | All   | 1620/2260 (72%) | 1385 (85%)   | 199 (12%)    | 36 (2%)    | 7           | 47 |

All 12 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

| Mol | Chain | Res | Type | Models (Total) |
|-----|-------|-----|------|----------------|
| 1   | B     | 70  | GLY  | 16             |
| 1   | B     | 69  | VAL  | 4              |
| 1   | B     | 19  | ILE  | 4              |
| 1   | B     | 42  | VAL  | 2              |
| 1   | B     | 18  | GLY  | 2              |
| 1   | B     | 37  | ILE  | 2              |
| 1   | B     | 84  | LYS  | 1              |
| 1   | B     | 16  | GLY  | 1              |
| 1   | B     | 14  | SER  | 1              |
| 1   | B     | 17  | LEU  | 1              |
| 1   | B     | 63  | VAL  | 1              |
| 1   | B     | 36  | GLY  | 1              |

### 6.3.2 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 PDB entries followed by that with respect to all NMR entries. 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   | B     | 66/90 (73%)     | 63±1 (95±2%) | 3±1 (5±2%) | 23          | 74 |
| All | All   | 1320/1800 (73%) | 1254 (95%)   | 66 (5%)    | 23          | 74 |

All 15 unique residues with a non-rotameric sidechain are listed below. They are sorted by the frequency of occurrence in the ensemble.

| Mol | Chain | Res | Type | Models (Total) |
|-----|-------|-----|------|----------------|
| 1   | B     | 5   | LEU  | 19             |
| 1   | B     | 10  | LEU  | 17             |
| 1   | B     | 41  | THR  | 11             |
| 1   | B     | 42  | VAL  | 4              |
| 1   | B     | 44  | GLU  | 2              |
| 1   | B     | 55  | GLN  | 2              |
| 1   | B     | 61  | VAL  | 2              |
| 1   | B     | 57  | ASN  | 2              |
| 1   | B     | 40  | LYS  | 1              |
| 1   | B     | 3   | MET  | 1              |
| 1   | B     | 11  | GLU  | 1              |
| 1   | B     | 24  | MET  | 1              |

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| Mol | Chain | Res | Type | Models (Total) |
|-----|-------|-----|------|----------------|
| 1   | B     | 53  | ARG  | 1              |
| 1   | B     | 12  | LYS  | 1              |
| 1   | B     | 82  | ASN  | 1              |

### 6.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

### 6.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

### 6.6 Ligand geometry [i](#)

There are no ligands in this entry.

### 6.7 Other polymers [i](#)

There are no such molecules in this entry.

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

There are no chain breaks in this entry.



## 7 Chemical shift validation

No chemical shift data were provided