



Full wwPDB NMR Structure Validation Report ⓘ

Mar 6, 2026 – 10:31 AM UTC

PDB ID : 1IBI / pdb_00001ibi
Title : QUAIL CYSTEINE AND GLYCINE-RICH PROTEIN, NMR, 15 MINIMIZED MODEL STRUCTURES
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Deposited on : 2001-03-28

This is a Full wwPDB NMR Structure Validation Report for a publicly released PDB entry.

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<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>.

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

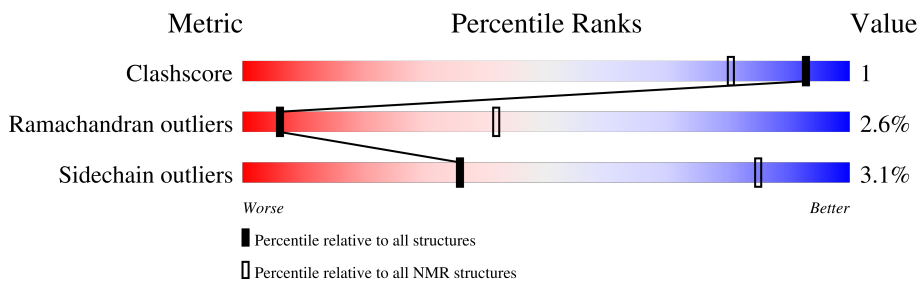
1 Overall quality at a glance

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 (#Entries)	NMR archive (#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 ≥ 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	A	113	

2 Ensemble composition and analysis

This entry contains 15 models. Model 7 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *lowest energy*.

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	A:118-A:175 (58)	1.03	7

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 3 clusters and 4 single-model clusters were found.

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

3 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 882 atoms, of which 434 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called CYSTEINE-RICH PROTEIN 2.

Mol	Chain	Residues	Atoms					Trace	
			Total	C	H	N	O		S
1	A	59	880	275	434	79	85	7	0

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	
2	A	2	Total	Zn
			2	2

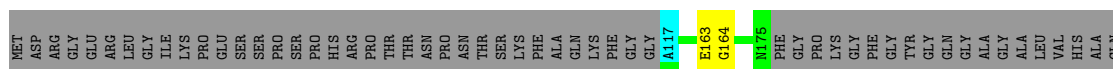
4 Residue-property plots [i](#)

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: CYSTEINE-RICH PROTEIN 2

Chain A:  50% 48%



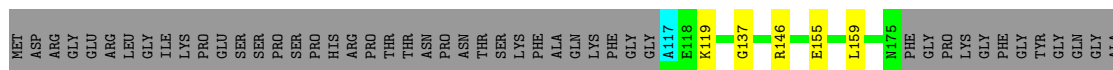
4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

- Molecule 1: CYSTEINE-RICH PROTEIN 2

Chain A:  47% 48%

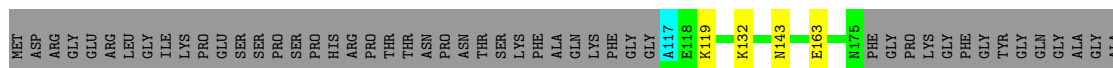


GLY
ALA
LEU
VAL
HIS
ALA
GLN

4.2.2 Score per residue for model 2

- Molecule 1: CYSTEINE-RICH PROTEIN 2

Chain A:  48% 48%



LEU
VAL
HIS
ALA
GLN

4.2.3 Score per residue for model 3

- Molecule 1: CYSTEINE-RICH PROTEIN 2

Chain A: 50% .. 48%

MET ASP ARG GLY GLU ARG LEU LEU LEU ILE ILE LYS PRO PRO GLU SER SER SER PRO PRO HIS HIS ARG PRO THR THR ASN ASN ASN THR SER SER PHE PHE ALA ALA LYS PHE GLY GLY A117 E118 T158 N175 PHE GLY PRO LYS LYS GLY PHE TYR GLY GLN GLY ALA ALA LEU VAL HIS GLN

4.2.4 Score per residue for model 4

- Molecule 1: CYSTEINE-RICH PROTEIN 2

Chain A: 47% . . 48%

MET ASP ARG GLY GLU ARG LEU LEU LEU ILE ILE LYS PRO PRO GLU SER SER SER PRO PRO HIS HIS ARG PRO THR THR ASN ASN ASN THR SER SER PHE PHE ALA ALA LYS PHE GLY GLY A117 K132 R146 K149 T158 C171 N175 PHE GLY PRO LYS LYS PHE PHE TYR GLY GLN GLY ALA

GLY
ALA
LEU
VAL
HIS
ALA
GLN

4.2.5 Score per residue for model 5

- Molecule 1: CYSTEINE-RICH PROTEIN 2

Chain A: 49% . . 48%

MET ASP ARG GLY GLU ARG LEU LEU LEU ILE ILE LYS PRO PRO GLU SER SER SER PRO PRO HIS HIS ARG PRO THR THR ASN ASN ASN THR SER SER PHE PHE ALA ALA LYS PHE GLY GLY A117 K132 T158 L159 N175 PHE GLY PRO LYS LYS PHE TYR GLY GLN GLY ALA ALA LEU VAL HIS

ALA
GLN

4.2.6 Score per residue for model 6

- Molecule 1: CYSTEINE-RICH PROTEIN 2

Chain A: 46% 5% . 48%

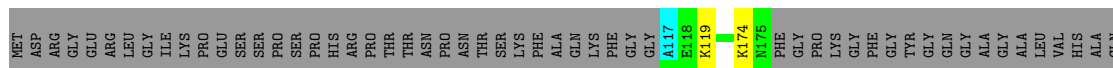
MET ASP ARG GLY GLU ARG LEU LEU LEU ILE ILE LYS PRO PRO GLU SER SER SER PRO PRO HIS HIS ARG PRO THR THR ASN ASN ASN THR SER SER PHE PHE ALA ALA LYS PHE GLY GLY A117 E118 K119 K132 G137 K149 L159 C168 N175 PHE GLY PRO LYS LYS PHE TYR GLY GLN GLY ALA ALA LEU VAL HIS GLN

GLY
ALA
LEU
VAL
HIS
ALA
GLN

4.2.7 Score per residue for model 7 (medoid)

- Molecule 1: CYSTEINE-RICH PROTEIN 2

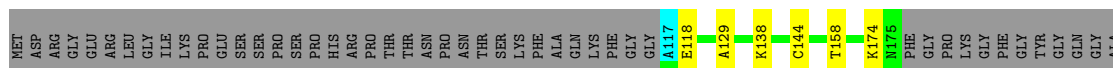
Chain A: 50% .. 48%



4.2.8 Score per residue for model 8

- Molecule 1: CYSTEINE-RICH PROTEIN 2

Chain A: 46% 5% . 48%

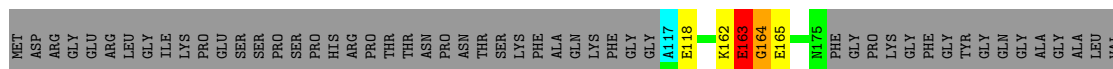


GLY
ALA
LEU
VAL
HIS
ALA
GLN

4.2.9 Score per residue for model 9

- Molecule 1: CYSTEINE-RICH PROTEIN 2

Chain A: 47% ... 48%

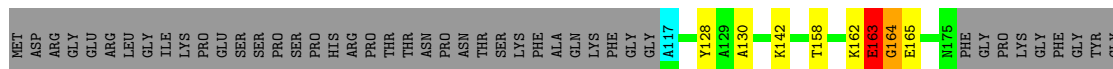


HIS
ALA
GLN

4.2.10 Score per residue for model 10

- Molecule 1: CYSTEINE-RICH PROTEIN 2

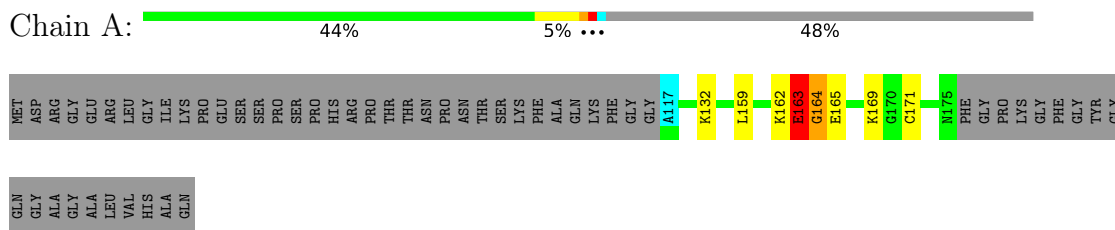
Chain A: 44% 5% ... 48%



GLN
GLY
ALA
GLY
ALA
LEU
VAL
HIS
ALA
GLN

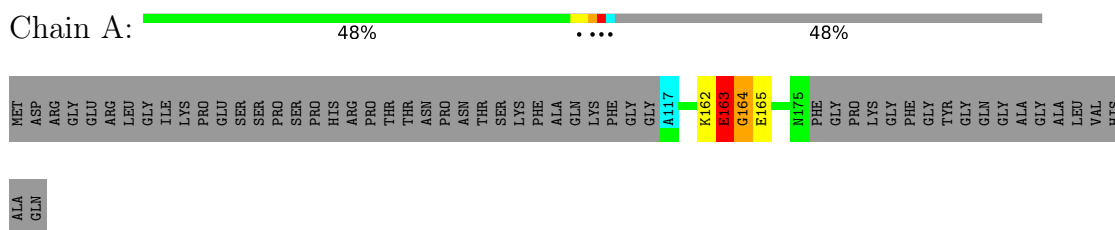
4.2.11 Score per residue for model 11

- Molecule 1: CYSTEINE-RICH PROTEIN 2



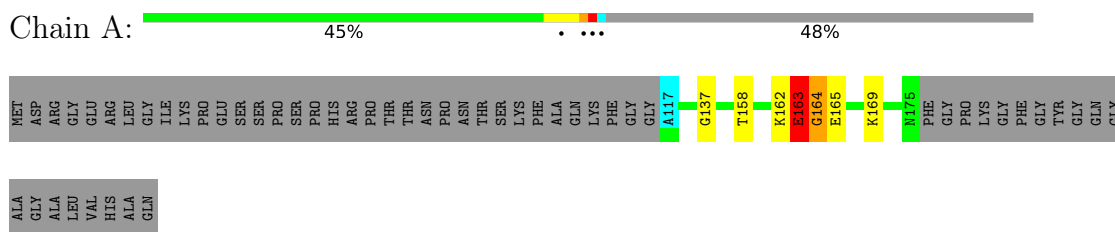
4.2.12 Score per residue for model 12

- Molecule 1: CYSTEINE-RICH PROTEIN 2



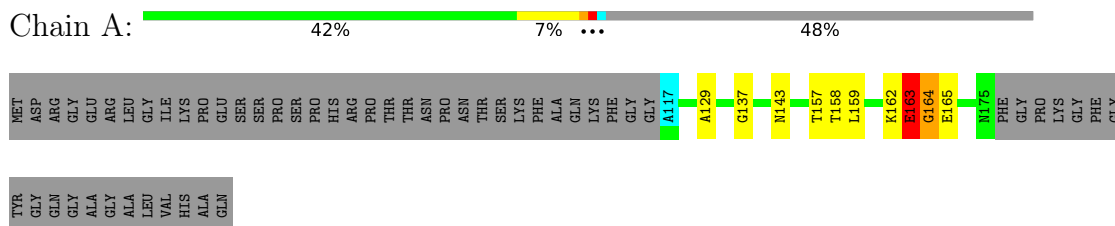
4.2.13 Score per residue for model 13

- Molecule 1: CYSTEINE-RICH PROTEIN 2



4.2.14 Score per residue for model 14

- Molecule 1: CYSTEINE-RICH PROTEIN 2



5 Refinement protocol and experimental data overview

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

Of the 100 calculated structures, 15 were deposited, based on the following criterion: *structures with the least restraint violations, structures with the lowest energy*.

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

Software name	Classification	Version
CNS	structure solution	1.0
X-PLOR	refinement	3.851

No chemical shift data was provided.

6 Model quality i

6.1 Standard geometry i

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

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	A	1.19±0.03	0±0/448 (0.1± 0.1%)	1.36±0.12	4±3/596 (0.6± 0.5%)
All	All	1.20	7/6720 (0.1%)	1.37	53/8940 (0.6%)

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	A	0.0±0.0	0.5±0.5
All	All	0	7

All unique bond outliers are listed below.

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
1	A	164	GLY	CA-C	5.75	1.59	1.51	15	7

All unique angle 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(°)	Ideal(°)	Models	
								Worst	Total
1	A	163	GLU	N-CA-CB	8.29	124.50	110.49	11	7
1	A	163	GLU	O-C-N	-7.89	112.09	122.59	13	7
1	A	128	TYR	N-CA-C	-7.04	104.56	113.43	10	1
1	A	162	LYS	CA-C-N	-6.79	108.58	121.54	11	7
1	A	162	LYS	C-N-CA	-6.79	108.58	121.54	11	7
1	A	158	THR	N-CA-C	-6.06	103.17	110.44	4	6
1	A	163	GLU	CA-C-N	-5.36	110.90	121.41	15	6
1	A	163	GLU	C-N-CA	-5.36	110.90	121.41	15	6

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	143	ASN	N-CA-C	-5.16	107.53	113.88	14	2
1	A	146	ARG	CA-C-N	5.08	128.09	120.87	1	1
1	A	146	ARG	C-N-CA	5.08	128.09	120.87	1	1
1	A	146	ARG	N-CA-CB	5.03	118.38	110.23	4	1
1	A	141	HIS	N-CA-CB	5.00	117.90	110.29	15	1

There are no chirality outliers.

All unique planar outliers are listed below.

Mol	Chain	Res	Type	Group	Models (Total)
1	A	163	GLU	Mainchain	7

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	A	441	429	430	1±1
2	A	2	0	0	0±0
All	All	6645	6435	6445	18

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 1.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:171:CYS:HG	2:A:196:ZN:ZN	0.92	0.64	4	3
1:A:144:CYS:HG	2:A:195:ZN:ZN	0.90	0.69	8	1
1:A:168:CYS:HG	2:A:196:ZN:ZN	0.87	0.59	6	1
1:A:159:LEU:HD23	1:A:159:LEU:H	0.46	1.70	5	3
1:A:163:GLU:C	1:A:165:GLU:H	0.45	2.20	12	3
1:A:163:GLU:O	1:A:165:GLU:N	0.44	2.51	13	7

6.3 Torsion angles [i](#)

6.3.1 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 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	A	57/113 (50%)	45±2 (79±3%)	11±2 (19±4%)	1±1 (3±3%)	6	42
All	All	855/1695 (50%)	672 (79%)	161 (19%)	22 (3%)	6	42

All 10 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	164	GLY	7
1	A	137	GLY	4
1	A	155	GLU	2
1	A	118	GLU	2
1	A	129	ALA	2
1	A	163	GLU	1
1	A	130	ALA	1
1	A	157	THR	1
1	A	158	THR	1
1	A	159	LEU	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	A	47/87 (54%)	46±1 (97±2%)	1±1 (3±2%)	36	85
All	All	705/1305 (54%)	683 (97%)	22 (3%)	36	85

All 10 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	A	132	LYS	5
1	A	119	LYS	4
1	A	169	LYS	3
1	A	149	LYS	2
1	A	174	LYS	2
1	A	142	LYS	2
1	A	159	LEU	1
1	A	118	GLU	1
1	A	138	LYS	1
1	A	128	TYR	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](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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