



# wwPDB NMR Structure Validation Summary Report ⓘ

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PDB ID : 1CMO / pdb\_00001cmo  
Title : IMMUNOGLOBULIN MOTIF DNA-RECOGNITION AND HET-  
ERODIMERIZATION FOR THE PEBP2/CBF RUNT-DOMAIN  
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This is a wwPDB NMR 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/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

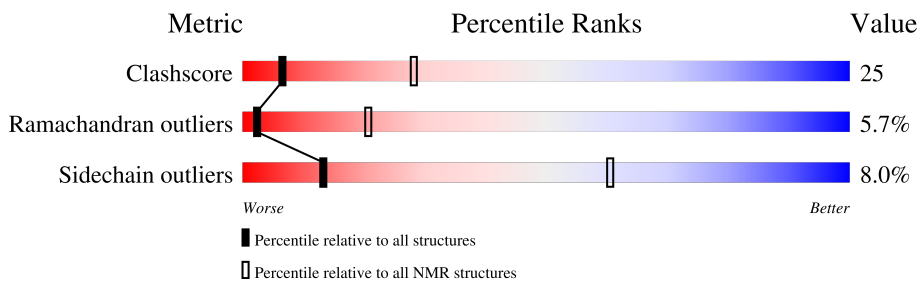
# 1 Overall quality at a glance

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 (#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  $\geq 3$ , 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	A	127	 51% 28% • 17%

## 2 Ensemble composition and analysis i

This entry contains 43 models. Model 21 is the overall representative, medoid model (most similar to other models).

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:62-A:166 (105)	0.61	21

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. No single-model clusters were found.

Cluster number	Models
1	3, 4, 6, 9, 11, 12, 13, 14, 16, 17, 18, 19, 20, 21, 22, 23, 24, 25, 26, 27, 28, 29, 30, 32, 33, 35, 37, 38, 39, 41
2	2, 8, 10, 31, 36, 42
3	7, 15, 34
4	40, 43
5	1, 5

### 3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 1977 atoms, of which 992 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called POLYOMAVIRUS ENHANCER BINDING PROTEIN 2.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	127	1977	619	992	182	181	3	0

There is a discrepancy between the modelled and reference sequences:

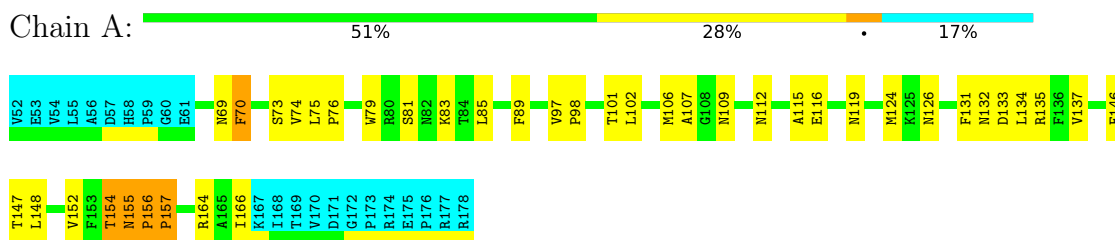
Chain	Residue	Modelled	Actual	Comment	Reference
A	81	SER	CYS	SEE REMARK 999	UNP Q01196

## 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.

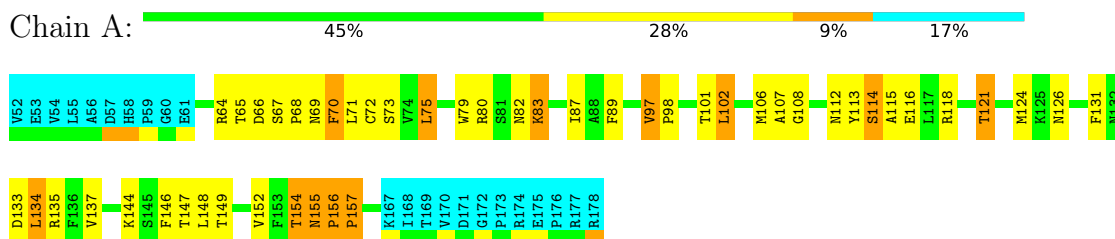
- Molecule 1: POLYOMAVIRUS ENHANCER BINDING PROTEIN 2



### 4.2 Residue scores for the representative (medoid) model from the NMR ensemble

The representative model is number 21. Colouring as in section 4.1 above.

- Molecule 1: POLYOMAVIRUS ENHANCER BINDING PROTEIN 2



## 5 Refinement protocol and experimental data overview

The models were refined using the following method: *HYBRID DISTANCE GEOMETRY/SIMULATED ANNEALING*.

Of the 300 calculated structures, 43 were deposited, based on the following criterion: *structures with the least restraint violations*.

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

Software name	Classification	Version
NMRPipe	structure solution	
PIPP	structure solution	
STAPP	structure solution	
X-PLOR	structure solution	
X-PLOR	refinement	3.843

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	A	1.55±0.01	0±0/830 ( 0.0± 0.1%)	1.24±0.01	1±1/1132 ( 0.0± 0.1%)
All	All	1.55	11/35690 ( 0.0%)	1.24	24/48676 ( 0.0%)

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	A	138	GLY	N-CA	7.84	1.50	1.44	20	3
1	A	108	GLY	N-CA	6.25	1.51	1.44	31	4
1	A	126	ASN	N-CA	5.76	1.50	1.46	40	2
1	A	95	GLY	N-CA	5.39	1.49	1.44	30	1
1	A	74	VAL	N-CA	5.14	1.52	1.46	43	1

5 of 7 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	153	PHE	CA-CB-CG	-5.71	108.09	113.80	30	7
1	A	163	HIS	CA-CB-CG	-5.46	108.34	113.80	27	6
1	A	84	THR	N-CA-C	-5.30	107.35	112.97	24	2
1	A	131	PHE	CA-CB-CG	-5.25	108.55	113.80	10	3
1	A	78	HIS	CA-CB-CG	-5.18	108.61	113.80	14	4

There are no chirality outliers.

There are no planarity outliers.

### 6.2 Too-close contacts [i](#)

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	812	819	819	41±8
All	All	34916	35217	35217	1760

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 25.

5 of 484 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:62:LEU:H	1:A:62:LEU:HD22	0.84	1.33	23	1
1:A:166:ILE:H	1:A:166:ILE:HD13	0.82	1.35	12	2
1:A:148:LEU:N	1:A:148:LEU:HD23	0.81	1.89	42	2
1:A:70:PHE:C	1:A:71:LEU:HD22	0.81	2.00	31	4
1:A:62:LEU:HD22	1:A:62:LEU:N	0.80	1.92	23	1

## 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	105/127 (83%)	82±7 (78±7%)	15±3 (14±2%)	6±2 (6±2%)	2	21
All	All	4424/5461 (81%)	3544 (80%)	627 (14%)	253 (6%)	2	21

5 of 19 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	156	PRO	43
1	A	157	PRO	41
1	A	83	LYS	27
1	A	133	ASP	18
1	A	86	PRO	18

### 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	89/108 (82%)	82±2 (92±3%)	7±2 (8±3%)	13	60
All	All	3827/4644 (82%)	3521 (92%)	306 (8%)	13	60

5 of 46 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	155	ASN	40
1	A	70	PHE	36
1	A	154	THR	28
1	A	148	LEU	20
1	A	79	TRP	16

### 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

There are no chain breaks in this entry.

## 7 Chemical shift validation

No chemical shift data were provided