



wwPDB NMR Structure Validation Summary Report ⓘ

Mar 5, 2026 – 08:47 AM UTC

PDB ID : 7RPM / pdb_00007rpm
BMRB ID : 30939
Title : Structures of the Intracellular Domain and Transmembrane Domain of the Human alpha7 Nicotinic Acetylcholine Receptors
Authors : Bondarenko, V.; Chen, Q.; Tang, P.
Deposited on : 2021-08-03

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

We welcome your comments at 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>.

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

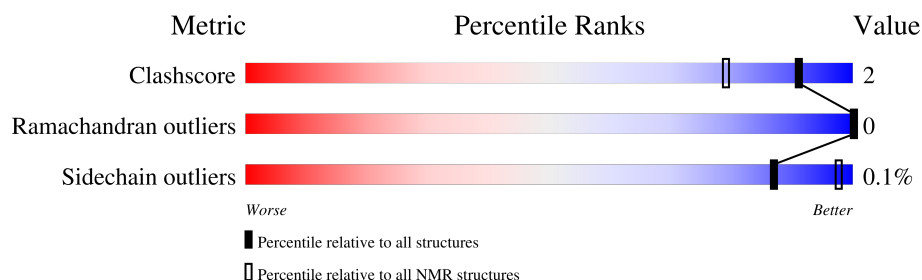
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

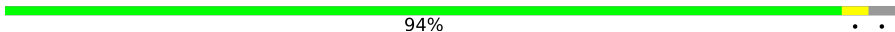
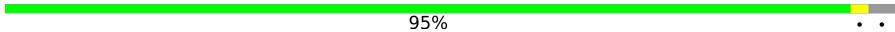
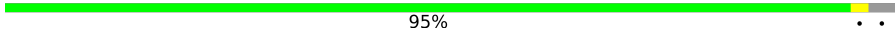
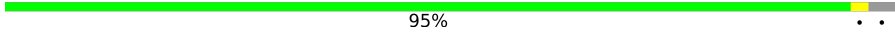
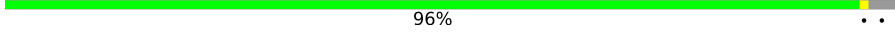
The overall completeness of chemical shifts assignment is 12%.

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	272	94% 
1	B	272	95% 
1	C	272	95% 
1	D	272	95% 
1	E	272	96% 

2 Ensemble composition and analysis

This entry contains 15 models. Model 9 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:209-A:472, B:209-B:472, C:209-C:472, D:209-D:472, E:209-E:472 (1320)	1.82	9

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

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

NmrClust was unable to cluster the ensemble.

Error message: Inconsistent models

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 20330 atoms, of which 10230 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Neuronal acetylcholine receptor subunit alpha-7.

Mol	Chain	Residues	Atoms						Trace
1	A	264	Total	C	H	N	O	S	0
			4066	1300	2046	340	359	21	
1	B	264	Total	C	H	N	O	S	0
			4066	1300	2046	340	359	21	
1	C	264	Total	C	H	N	O	S	0
			4066	1300	2046	340	359	21	
1	D	264	Total	C	H	N	O	S	0
			4066	1300	2046	340	359	21	
1	E	264	Total	C	H	N	O	S	0
			4066	1300	2046	340	359	21	

There are 50 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	203	SER	-	expression tag	UNP P36544
A	204	ASN	-	expression tag	UNP P36544
A	205	ALA	-	expression tag	UNP P36544
A	206	GLU	-	expression tag	UNP P36544
A	207	GLU	-	expression tag	UNP P36544
A	208	GLU	-	expression tag	UNP P36544
A	263	SER	ALA	engineered mutation	UNP P36544
A	268	SER	VAL	engineered mutation	UNP P36544
A	270	SER	LEU	engineered mutation	UNP P36544
A	474	GLU	ALA	engineered mutation	UNP P36544
B	203	SER	-	expression tag	UNP P36544
B	204	ASN	-	expression tag	UNP P36544
B	205	ALA	-	expression tag	UNP P36544
B	206	GLU	-	expression tag	UNP P36544
B	207	GLU	-	expression tag	UNP P36544
B	208	GLU	-	expression tag	UNP P36544
B	263	SER	ALA	engineered mutation	UNP P36544
B	268	SER	VAL	engineered mutation	UNP P36544
B	270	SER	LEU	engineered mutation	UNP P36544
B	474	GLU	ALA	engineered mutation	UNP P36544
C	203	SER	-	expression tag	UNP P36544
C	204	ASN	-	expression tag	UNP P36544
C	205	ALA	-	expression tag	UNP P36544

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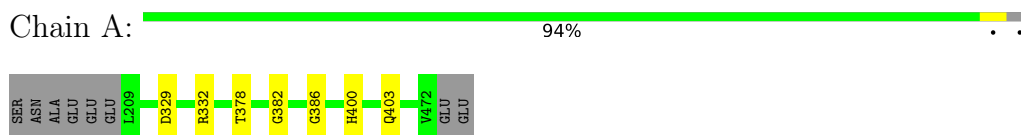
Chain	Residue	Modelled	Actual	Comment	Reference
C	206	GLU	-	expression tag	UNP P36544
C	207	GLU	-	expression tag	UNP P36544
C	208	GLU	-	expression tag	UNP P36544
C	263	SER	ALA	engineered mutation	UNP P36544
C	268	SER	VAL	engineered mutation	UNP P36544
C	270	SER	LEU	engineered mutation	UNP P36544
C	474	GLU	ALA	engineered mutation	UNP P36544
D	203	SER	-	expression tag	UNP P36544
D	204	ASN	-	expression tag	UNP P36544
D	205	ALA	-	expression tag	UNP P36544
D	206	GLU	-	expression tag	UNP P36544
D	207	GLU	-	expression tag	UNP P36544
D	208	GLU	-	expression tag	UNP P36544
D	263	SER	ALA	engineered mutation	UNP P36544
D	268	SER	VAL	engineered mutation	UNP P36544
D	270	SER	LEU	engineered mutation	UNP P36544
D	474	GLU	ALA	engineered mutation	UNP P36544
E	203	SER	-	expression tag	UNP P36544
E	204	ASN	-	expression tag	UNP P36544
E	205	ALA	-	expression tag	UNP P36544
E	206	GLU	-	expression tag	UNP P36544
E	207	GLU	-	expression tag	UNP P36544
E	208	GLU	-	expression tag	UNP P36544
E	263	SER	ALA	engineered mutation	UNP P36544
E	268	SER	VAL	engineered mutation	UNP P36544
E	270	SER	LEU	engineered mutation	UNP P36544
E	474	GLU	ALA	engineered mutation	UNP P36544

4 Residue-property plots [i](#)

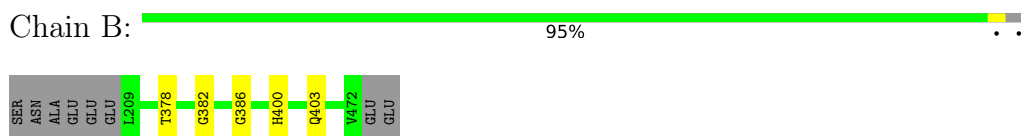
4.1 Average score per residue in the NMR ensemble

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.

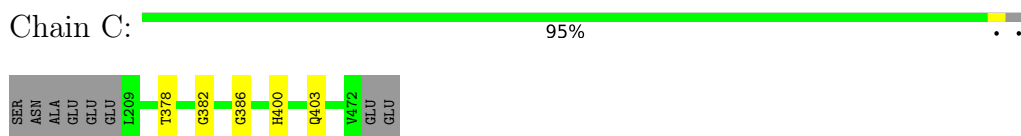
- Molecule 1: Neuronal acetylcholine receptor subunit alpha-7



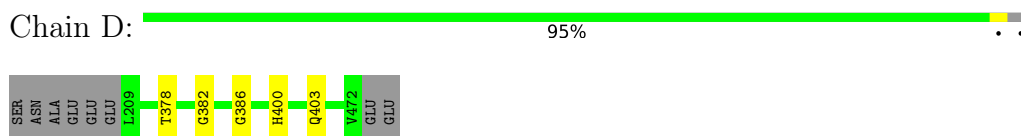
- Molecule 1: Neuronal acetylcholine receptor subunit alpha-7



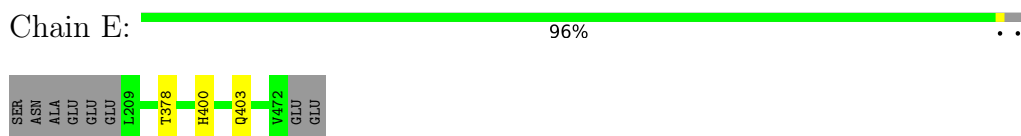
- Molecule 1: Neuronal acetylcholine receptor subunit alpha-7



- Molecule 1: Neuronal acetylcholine receptor subunit alpha-7



- Molecule 1: Neuronal acetylcholine receptor subunit alpha-7



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

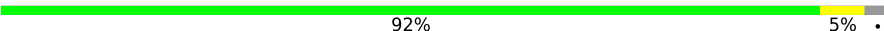
The representative model is number 9. Colouring as in section 4.1 above.

- Molecule 1: Neuronal acetylcholine receptor subunit alpha-7

Chain A:  91% 6% .



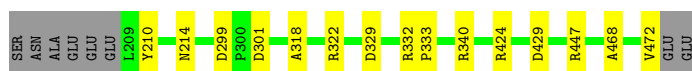
- Molecule 1: Neuronal acetylcholine receptor subunit alpha-7

Chain B:  92% 5% .

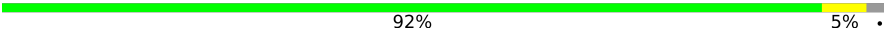


- Molecule 1: Neuronal acetylcholine receptor subunit alpha-7

Chain C:  92% 6% .



- Molecule 1: Neuronal acetylcholine receptor subunit alpha-7

Chain D:  92% 5% .



- Molecule 1: Neuronal acetylcholine receptor subunit alpha-7

Chain E:  92% 5% .



5 Refinement protocol and experimental data overview

The models were refined using the following method: *na, na*.

Of the 1000 calculated structures, 15 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
CS-ROSETTA	structure calculation	3.17
Rosetta	structure calculation	3.7
Rosetta	refinement	3.7
PHENIX	refinement	1.19

The following table shows chemical shift validation statistics as aggregates over all chemical shift files. Detailed validation can be found in section 7 of this report.

Chemical shift file(s)	working_cs.cif
Number of chemical shift lists	1
Total number of shifts	2113
Number of shifts mapped to atoms	2113
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	12%

6 Model quality [i](#)

6.1 Standard geometry [i](#)

There are no covalent bond-length or bond-angle outliers.

There are no bond-length outliers.

There are no bond-angle outliers.

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	2020	2046	2054	8±3
1	B	2020	2046	2054	8±3
1	C	2020	2046	2054	8±3
1	D	2020	2046	2054	8±3
1	E	2020	2046	2054	8±3
All	All	151500	153450	154050	490

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 2.

5 of 246 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:406:GLU:O	1:E:400:HIS:NE2	0.81	2.14	14	5
1:B:400:HIS:NE2	1:C:406:GLU:O	0.81	2.14	14	5
1:D:400:HIS:NE2	1:E:406:GLU:O	0.79	2.14	14	5
1:A:400:HIS:NE2	1:B:406:GLU:O	0.79	2.16	6	5
1:C:400:HIS:NE2	1:D:406:GLU:O	0.78	2.14	14	5

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	A	262/272 (96%)	257±2 (98±1%)	5±2 (2±1%)	0±0 (0±0%)	100	100
1	B	262/272 (96%)	257±2 (98±1%)	5±2 (2±1%)	0±0 (0±0%)	100	100
1	C	262/272 (96%)	257±2 (98±1%)	5±2 (2±1%)	0±0 (0±0%)	100	100
1	D	262/272 (96%)	257±2 (98±1%)	5±2 (2±1%)	0±0 (0±0%)	100	100
1	E	262/272 (96%)	257±2 (98±1%)	5±2 (2±1%)	0±0 (0±0%)	100	100
All	All	19650/20400 (96%)	19259 (98%)	391 (2%)	0 (0%)	100	100

There are no Ramachandran outliers.

6.3.2 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 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	227/234 (97%)	227±0 (100±0%)	0±0 (0±0%)	87	98
1	B	227/234 (97%)	227±0 (100±0%)	0±0 (0±0%)	87	98
1	C	227/234 (97%)	227±0 (100±0%)	0±0 (0±0%)	87	98
1	D	227/234 (97%)	227±0 (100±0%)	0±0 (0±0%)	87	98
1	E	227/234 (97%)	227±0 (100±0%)	0±0 (0±0%)	87	98
All	All	17025/17550 (97%)	17005 (100%)	20 (0%)	87	98

5 of 20 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	423	ASN	1
1	B	423	ASN	1

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Mol	Chain	Res	Type	Models (Total)
1	C	423	ASN	1
1	D	423	ASN	1
1	E	423	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 [i](#)

The completeness of assignment taking into account all chemical shift lists is 12% for the well-defined parts and 12% for the entire structure.

7.1 Chemical shift list 1

File name: `working_cs.cif`

Chemical shift list name: *starch_output*

7.1.1 Bookkeeping [i](#)

The following table shows the results of parsing the chemical shift list and reports the number of nuclei with statistically unusual chemical shifts.

Total number of shifts	2113
Number of shifts mapped to atoms	2113
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

7.1.2 Chemical shift referencing [i](#)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	259	-0.51 ± 0.09	Should be checked
$^{13}\text{C}_\beta$	170	0.24 ± 0.02	None needed (< 0.5 ppm)
$^{13}\text{C}'$	200	-0.09 ± 0.08	None needed (< 0.5 ppm)
^{15}N	231	0.46 ± 0.13	None needed (< 0.5 ppm)

7.1.3 Completeness of resonance assignments [i](#)

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 12%, i.e. 2113 atoms were assigned a chemical shift out of a possible 17800. 0 out of 285 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	1155/6505 (18%)	465/2635 (18%)	459/2640 (17%)	231/1230 (19%)
Sidechain	910/10045 (9%)	618/6665 (9%)	288/3100 (9%)	4/280 (1%)

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	Total	^1H	^{13}C	^{15}N
Aromatic	48/1250 (4%)	27/650 (4%)	17/580 (3%)	4/20 (20%)
Overall	2113/17800 (12%)	1110/9950 (11%)	764/6320 (12%)	239/1530 (16%)

7.1.4 Statistically unusual chemical shifts [i](#)

There are no statistically unusual chemical shifts.

7.1.5 Random Coil Index (RCI) plots [i](#)

The image below reports *random coil index* values for the protein chains in the structure. The height of each bar gives a probability of a given residue to be disordered, as predicted from the available chemical shifts and the amino acid sequence. A value above 0.2 is an indication of significant predicted disorder. The colour of the bar shows whether the residue is in the well-defined core (black) or in the ill-defined residue ranges (cyan), as described in section 2 on ensemble composition. If well-defined core and ill-defined regions are not identified then it is shown as gray bars.

Random coil index (RCI) for chain A:

