



Full wwPDB NMR Structure Validation Report ⓘ

Mar 10, 2026 – 10:54 AM UTC

PDB ID : 1YT6 / pdb_00001yt6
Title : NMR structure of peptide SD
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Deposited on : 2005-02-10

This is a Full wwPDB NMR Structure Validation 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

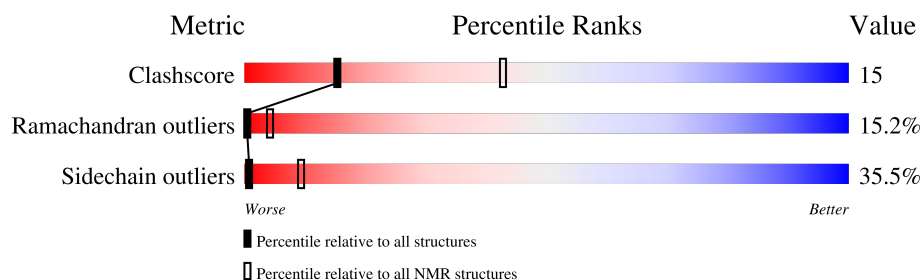
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment 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	10	50% 40% 10%

2 Ensemble composition and analysis ⓘ

This entry contains 50 models.

Cyrange was unable to find well-defined residues.

Error message: The number of core atoms (5) was below the domain threshold value (8).

The models can be grouped into 6 clusters and 3 single-model clusters were found.

Cluster number	Models
1	7, 8, 9, 14, 17, 25, 28, 30, 31, 37, 38, 39, 44, 48, 49, 50
2	1, 6, 11, 12, 13, 18, 21, 22, 26, 29, 33, 40, 41, 43, 47
3	2, 4, 10, 35, 36, 42
4	5, 15, 20, 23
5	24, 27, 32, 45
6	34, 46
Single-model clusters	3; 16; 19

3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 134 atoms, of which 62 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called peptide SD.

Mol	Chain	Residues	Atoms						Trace
1	A	10	Total	C	H	N	O	S	
			134	45	62	11	14	2	0

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: peptide SD



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: peptide SD



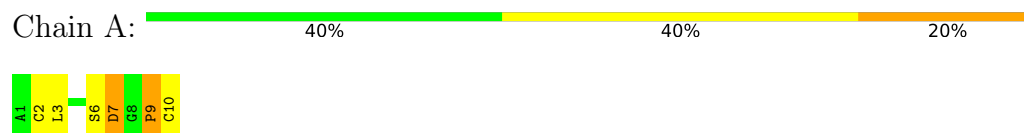
4.2.2 Score per residue for model 2

- Molecule 1: peptide SD



4.2.3 Score per residue for model 3

- Molecule 1: peptide SD



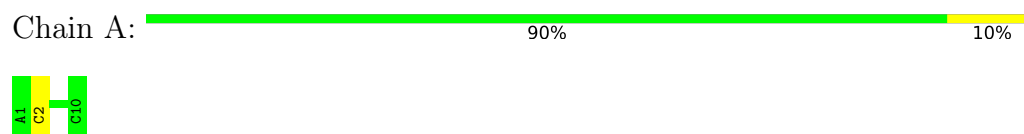
4.2.4 Score per residue for model 4

- Molecule 1: peptide SD



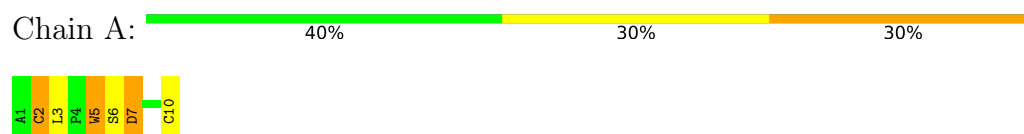
4.2.5 Score per residue for model 5

- Molecule 1: peptide SD



4.2.6 Score per residue for model 6

- Molecule 1: peptide SD



4.2.7 Score per residue for model 7

- Molecule 1: peptide SD



4.2.8 Score per residue for model 8

- Molecule 1: peptide SD



4.2.9 Score per residue for model 9

- Molecule 1: peptide SD



4.2.10 Score per residue for model 10

- Molecule 1: peptide SD



4.2.11 Score per residue for model 11

- Molecule 1: peptide SD



4.2.12 Score per residue for model 12

- Molecule 1: peptide SD



4.2.13 Score per residue for model 13

- Molecule 1: peptide SD



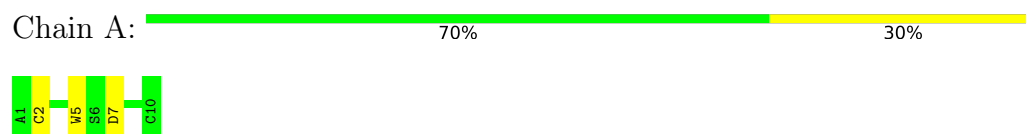
4.2.14 Score per residue for model 14

- Molecule 1: peptide SD



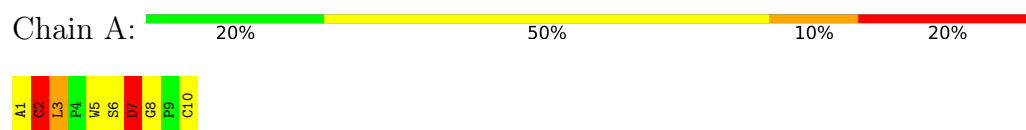
4.2.15 Score per residue for model 15

- Molecule 1: peptide SD



4.2.16 Score per residue for model 16

- Molecule 1: peptide SD



4.2.17 Score per residue for model 17

- Molecule 1: peptide SD



4.2.18 Score per residue for model 18

- Molecule 1: peptide SD



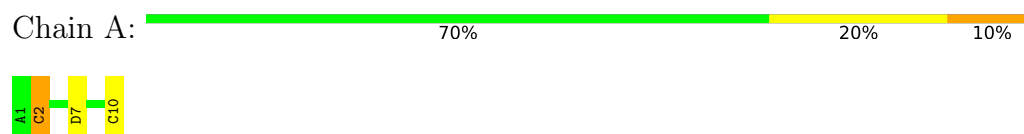
4.2.19 Score per residue for model 19

- Molecule 1: peptide SD



4.2.20 Score per residue for model 20

- Molecule 1: peptide SD



4.2.21 Score per residue for model 21

- Molecule 1: peptide SD



4.2.22 Score per residue for model 22

- Molecule 1: peptide SD



4.2.23 Score per residue for model 23

- Molecule 1: peptide SD



4.2.24 Score per residue for model 24

- Molecule 1: peptide SD



4.2.25 Score per residue for model 25

- Molecule 1: peptide SD



4.2.26 Score per residue for model 26

- Molecule 1: peptide SD



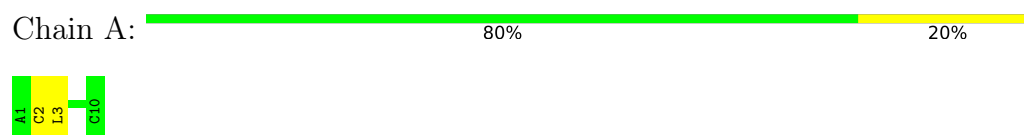
4.2.27 Score per residue for model 27

- Molecule 1: peptide SD



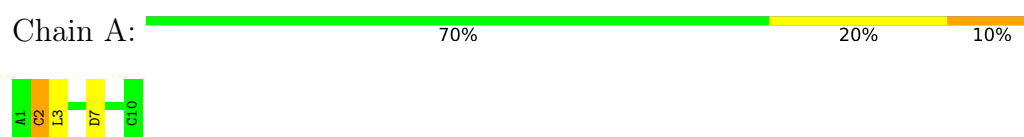
4.2.28 Score per residue for model 28

- Molecule 1: peptide SD



4.2.29 Score per residue for model 29

- Molecule 1: peptide SD



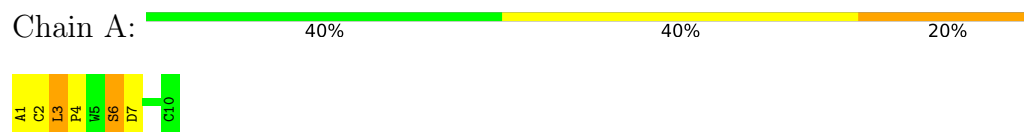
4.2.30 Score per residue for model 30

- Molecule 1: peptide SD



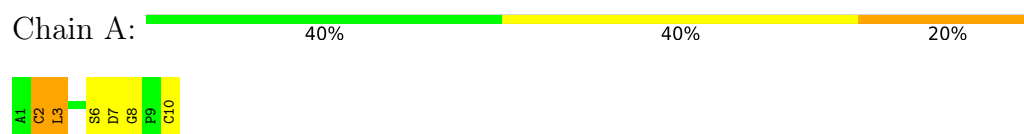
4.2.31 Score per residue for model 31

- Molecule 1: peptide SD



4.2.32 Score per residue for model 32

- Molecule 1: peptide SD



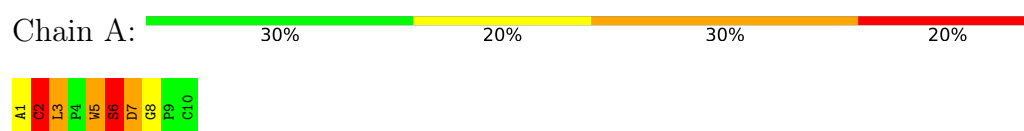
4.2.33 Score per residue for model 33

- Molecule 1: peptide SD



4.2.34 Score per residue for model 34

- Molecule 1: peptide SD



4.2.35 Score per residue for model 35

- Molecule 1: peptide SD



4.2.36 Score per residue for model 36

- Molecule 1: peptide SD



4.2.37 Score per residue for model 37

- Molecule 1: peptide SD



4.2.38 Score per residue for model 38

- Molecule 1: peptide SD



4.2.39 Score per residue for model 39

- Molecule 1: peptide SD



4.2.40 Score per residue for model 40

- Molecule 1: peptide SD



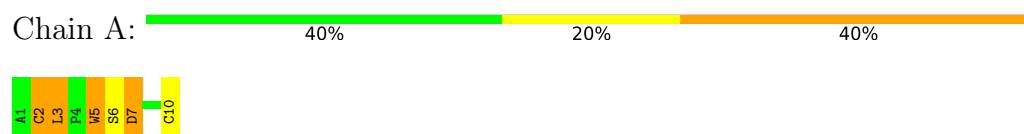
4.2.41 Score per residue for model 41

- Molecule 1: peptide SD



4.2.42 Score per residue for model 42

- Molecule 1: peptide SD



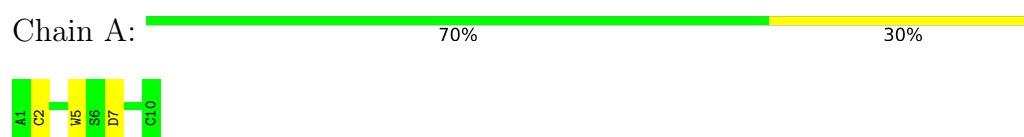
4.2.43 Score per residue for model 43

- Molecule 1: peptide SD



4.2.44 Score per residue for model 44

- Molecule 1: peptide SD



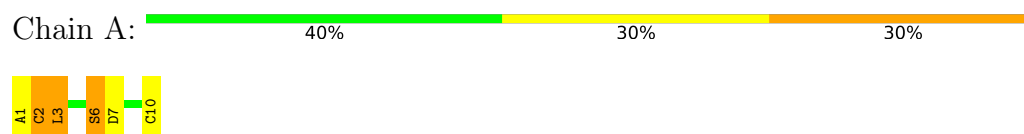
4.2.45 Score per residue for model 45

- Molecule 1: peptide SD



4.2.46 Score per residue for model 46

- Molecule 1: peptide SD



4.2.47 Score per residue for model 47

- Molecule 1: peptide SD



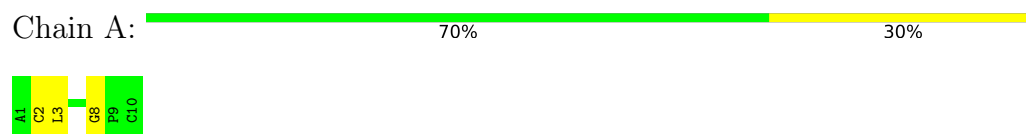
4.2.48 Score per residue for model 48

- Molecule 1: peptide SD



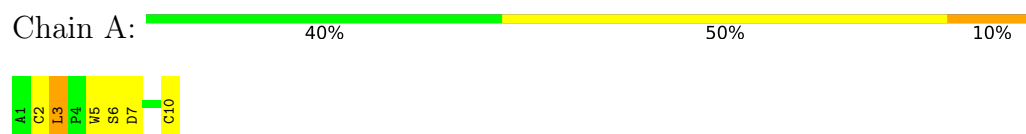
4.2.49 Score per residue for model 49

- Molecule 1: peptide SD



4.2.50 Score per residue for model 50

- Molecule 1: peptide SD



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, 50 were deposited, based on the following criterion: *all calculated structures submitted*.

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

Software name	Classification	Version
CNS	refinement	1.0

No chemical shift data was provided.

6 Model quality ⓘ

6.1 Standard geometry ⓘ

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 ⓘ

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	72	62	62	2±2
All	All	3600	3100	3100	99

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 15.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:3:LEU:HD12	1:A:5:TRP:CZ3	0.84	2.07	10	2
1:A:3:LEU:HD12	1:A:6:SER:CB	0.79	2.08	45	2
1:A:3:LEU:HD12	1:A:6:SER:HB2	0.79	1.54	45	4
1:A:2:CYS:O	1:A:3:LEU:HD23	0.78	1.79	30	4
1:A:2:CYS:C	1:A:3:LEU:HD23	0.70	2.12	24	5
1:A:1:ALA:HB3	1:A:7:ASP:HA	0.68	1.65	14	2
1:A:3:LEU:HD13	1:A:5:TRP:CZ3	0.68	2.23	6	5
1:A:3:LEU:HD23	1:A:4:PRO:HD2	0.67	1.65	23	3
1:A:2:CYS:C	1:A:3:LEU:HD12	0.67	2.14	43	1
1:A:3:LEU:HD12	1:A:5:TRP:CH2	0.63	2.29	42	2
1:A:3:LEU:HD23	1:A:3:LEU:N	0.63	2.08	16	2
1:A:3:LEU:HD12	1:A:6:SER:HB3	0.63	1.70	19	3
1:A:3:LEU:HD22	1:A:6:SER:CB	0.62	2.24	50	2
1:A:3:LEU:HB2	1:A:7:ASP:CB	0.60	2.25	45	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:3:LEU:HD23	1:A:7:ASP:O	0.58	1.98	31	3
1:A:3:LEU:N	1:A:3:LEU:HD12	0.56	2.16	21	2
1:A:3:LEU:HD13	1:A:7:ASP:O	0.56	2.01	35	2
1:A:3:LEU:HD22	1:A:5:TRP:CZ3	0.55	2.36	48	1
1:A:3:LEU:HD12	1:A:6:SER:O	0.55	2.01	47	1
1:A:3:LEU:HD22	1:A:6:SER:HB2	0.55	1.78	50	1
1:A:6:SER:O	1:A:8:GLY:N	0.55	2.40	27	4
1:A:6:SER:O	1:A:7:ASP:C	0.55	2.50	43	5
1:A:3:LEU:HD11	1:A:9:PRO:O	0.53	2.03	3	1
1:A:3:LEU:HD21	1:A:9:PRO:CG	0.51	2.35	30	1
1:A:7:ASP:C	1:A:9:PRO:HD3	0.51	2.30	30	1
1:A:3:LEU:HD13	1:A:5:TRP:HZ3	0.51	1.65	18	2
1:A:3:LEU:HD11	1:A:9:PRO:HG3	0.50	1.84	30	1
1:A:1:ALA:HB1	1:A:6:SER:HA	0.50	1.82	30	1
1:A:2:CYS:N	1:A:7:ASP:O	0.49	2.46	30	2
1:A:3:LEU:HD22	1:A:7:ASP:CB	0.49	2.38	3	1
1:A:3:LEU:HD13	1:A:5:TRP:CH2	0.48	2.43	6	1
1:A:3:LEU:HD12	1:A:7:ASP:C	0.48	2.34	8	1
1:A:3:LEU:N	1:A:3:LEU:CD2	0.46	2.79	16	2
1:A:1:ALA:HB1	1:A:7:ASP:CB	0.46	2.40	33	1
1:A:3:LEU:HD13	1:A:6:SER:HB2	0.46	1.86	23	1
1:A:3:LEU:HD12	1:A:6:SER:OG	0.46	2.11	34	1
1:A:3:LEU:CD1	1:A:5:TRP:CZ3	0.45	3.00	34	1
1:A:3:LEU:HB2	1:A:7:ASP:C	0.45	2.36	34	1
1:A:3:LEU:HD22	1:A:6:SER:HB3	0.45	1.89	50	1
1:A:3:LEU:CD2	1:A:5:TRP:CZ3	0.45	3.00	50	1
1:A:3:LEU:HD13	1:A:4:PRO:CD	0.45	2.41	31	1
1:A:3:LEU:HD23	1:A:4:PRO:CD	0.45	2.41	41	1
1:A:3:LEU:HD23	1:A:6:SER:HB2	0.44	1.89	39	1
1:A:2:CYS:HB3	1:A:3:LEU:HD12	0.44	1.89	43	1
1:A:3:LEU:HB2	1:A:6:SER:O	0.44	2.12	43	1
1:A:6:SER:O	1:A:7:ASP:O	0.44	2.36	6	1
1:A:1:ALA:O	1:A:2:CYS:C	0.44	2.61	26	9
1:A:3:LEU:HG	1:A:7:ASP:HB2	0.43	1.90	16	1
1:A:3:LEU:HD12	1:A:8:GLY:HA3	0.43	1.89	37	1
1:A:3:LEU:N	1:A:3:LEU:CD1	0.42	2.82	27	1
1:A:1:ALA:HB1	1:A:6:SER:CA	0.42	2.45	31	1
1:A:3:LEU:CD1	1:A:3:LEU:N	0.41	2.83	28	1
1:A:3:LEU:CD1	1:A:7:ASP:O	0.40	2.69	2	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	A	8/10 (80%)	4±1 (50±13%)	3±1 (34±13%)	1±1 (15±12%)	0	4
All	All	400/500 (80%)	202 (50%)	137 (34%)	61 (15%)	0	4

All 5 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	7	ASP	19
1	A	6	SER	18
1	A	2	CYS	12
1	A	8	GLY	10
1	A	9	PRO	2

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	8/8 (100%)	5±1 (64±12%)	3±1 (36±12%)	1	9
All	All	400/400 (100%)	258 (64%)	142 (36%)	1	9

All 6 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	2	CYS	49
1	A	5	TRP	28
1	A	7	ASP	23
1	A	3	LEU	22
1	A	10	CYS	14

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Mol	Chain	Res	Type	Models (Total)
1	A	6	SER	6

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