



Full wwPDB NMR Structure Validation Report ⓘ

Mar 9, 2026 – 05:07 PM UTC

PDB ID : 6MI5 / pdb_00006mi5
BMRB ID : 30515
Title : NMR solution structure of lanmodulin (LanM) complexed with yttrium(III) ions
Authors : Cook, E.C.; Featherson, E.R.; Showalter, S.A.; Cotruvo Jr., J.A.
Deposited on : 2018-09-19

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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

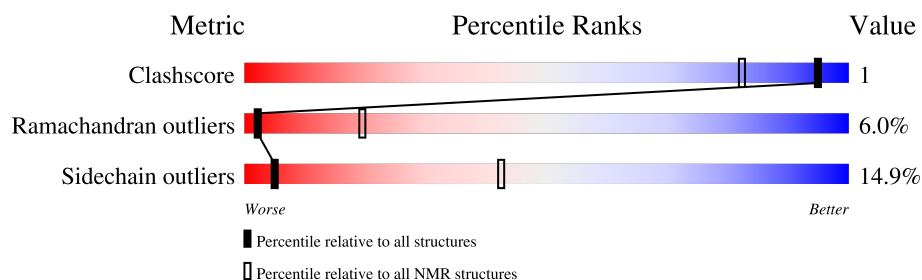
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 72%.

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	X	117	

2 Ensemble composition and analysis

This entry contains 12 models. Model 2 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *minimized average structure*.

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	X:28-X:134 (107)	1.02	2

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 2 clusters and 4 single-model clusters were found.

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

3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1739 atoms, of which 854 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Lanmodulin.

Mol	Chain	Residues	Atoms					Trace
1	X	117	Total	C	H	N	O	0
			1736	543	854	155	184	

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
X	134	HIS	-	expression tag	UNP C5B164
X	135	HIS	-	expression tag	UNP C5B164
X	136	HIS	-	expression tag	UNP C5B164
X	137	HIS	-	expression tag	UNP C5B164
X	138	HIS	-	expression tag	UNP C5B164
X	139	HIS	-	expression tag	UNP C5B164

- Molecule 2 is YTTRIUM (III) ION (CCD ID: YT3) (formula: Y).

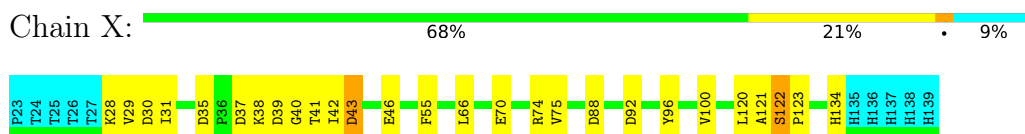
Mol	Chain	Residues	Atoms	
2	X	3	Total	Y
			3	3

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: Lanmodulin

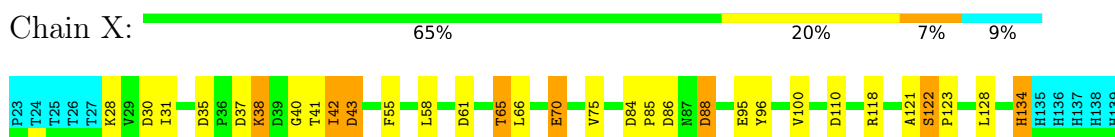


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

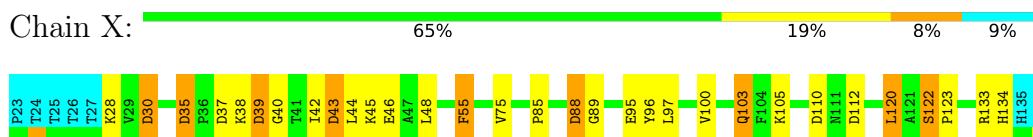
4.2.1 Score per residue for model 1

- Molecule 1: Lanmodulin



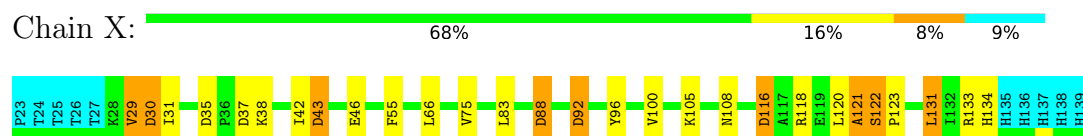
4.2.2 Score per residue for model 2 (medoid)

- Molecule 1: Lanmodulin



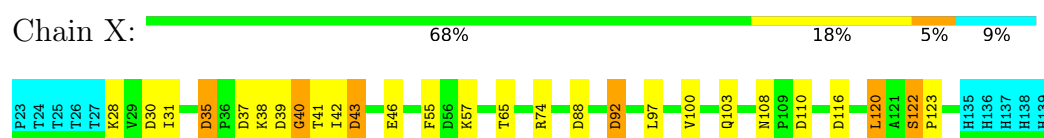
4.2.3 Score per residue for model 3

- Molecule 1: Lanmodulin



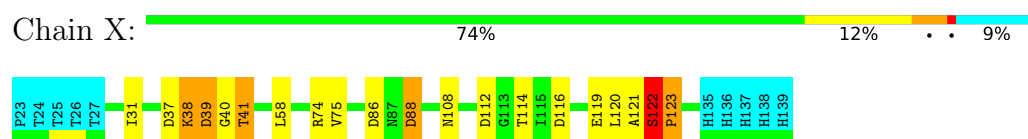
4.2.4 Score per residue for model 4

- Molecule 1: Lanmodulin



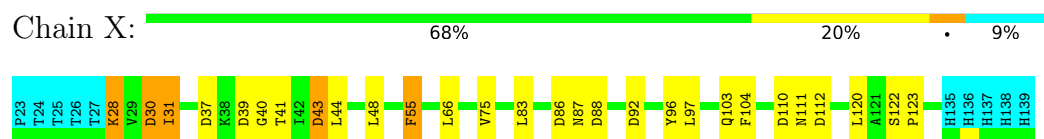
4.2.5 Score per residue for model 5

- Molecule 1: Lanmodulin



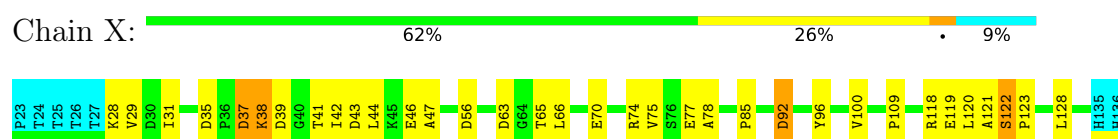
4.2.6 Score per residue for model 6

- Molecule 1: Lanmodulin



4.2.7 Score per residue for model 7

- Molecule 1: Lanmodulin

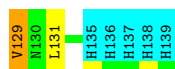
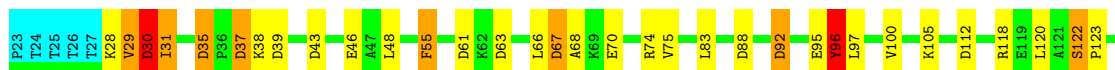




4.2.8 Score per residue for model 8

- Molecule 1: Lanmodulin

Chain X: 62% 21% 8% 9%



4.2.9 Score per residue for model 9

- Molecule 1: Lanmodulin

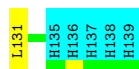
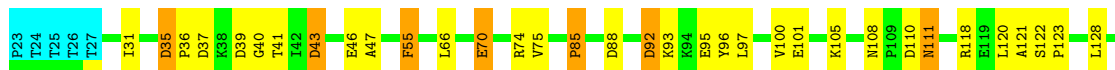
Chain X: 71% 15% 6% 9%



4.2.10 Score per residue for model 10

- Molecule 1: Lanmodulin

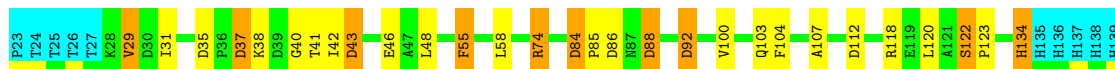
Chain X: 62% 24% 6% 9%



4.2.11 Score per residue for model 11

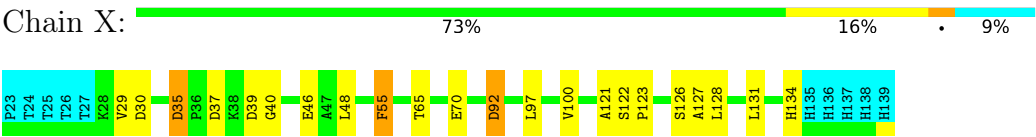
- Molecule 1: Lanmodulin

Chain X: 67% 16% 9% 9%



4.2.12 Score per residue for model 12

● Molecule 1: Lanmodulin



5 Refinement protocol and experimental data overview

The models were refined using the following method: *molecular dynamics*.

Of the 12 calculated structures, 12 were deposited, based on the following criterion: *structures with acceptable covalent geometry*.

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

Software name	Classification	Version
Amber	refinement	
Amber	structure calculation	14

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	998
Number of shifts mapped to atoms	998
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	72%

6 Model quality

6.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: YT3

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	X	0.88±0.02	0±0/806 (0.0± 0.0%)	1.73±0.04	11±3/1090 (1.0± 0.3%)
All	All	0.88	0/9672 (0.0%)	1.73	129/13080 (1.0%)

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	X	0.0±0.0	6.3±1.9
All	All	0	76

There are no bond-length outliers.

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	X	43	ASP	CA-CB-CG	9.22	121.82	112.60	9	9
1	X	88	ASP	CA-CB-CG	8.60	121.20	112.60	1	5
1	X	92	ASP	CA-CB-CG	7.92	120.52	112.60	7	7
1	X	37	ASP	CA-CB-CG	7.68	120.28	112.60	4	9
1	X	116	ASP	CA-CB-CG	7.45	120.05	112.60	3	2
1	X	55	PHE	CA-CB-CG	7.36	121.16	113.80	9	6
1	X	92	ASP	OD1-CG-OD2	-6.92	106.28	122.90	8	5
1	X	111	ASN	CA-CB-CG	6.79	119.39	112.60	6	1
1	X	85	PRO	CA-C-N	6.56	133.51	121.70	10	1
1	X	85	PRO	C-N-CA	6.56	133.51	121.70	10	1
1	X	86	ASP	CA-CB-CG	6.55	119.15	112.60	5	3
1	X	46	GLU	OE1-CD-OE2	-6.54	107.22	122.90	7	9

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	X	108	ASN	CA-CB-CG	6.17	118.77	112.60	5	1
1	X	67	ASP	CA-CB-CG	6.12	118.72	112.60	8	1
1	X	47	ALA	CA-C-N	6.06	128.31	120.44	10	2
1	X	47	ALA	C-N-CA	6.06	128.31	120.44	10	2
1	X	37	ASP	OD1-CG-OD2	-5.95	108.61	122.90	9	10
1	X	39	ASP	OD1-CG-OD2	-5.91	108.72	122.90	8	6
1	X	29	VAL	N-CA-C	5.89	116.67	110.72	12	1
1	X	111	ASN	N-CA-C	5.84	123.24	110.80	10	1
1	X	133	ARG	N-CA-C	5.82	118.37	111.33	9	1
1	X	89	GLY	CA-C-N	5.78	131.35	121.87	2	1
1	X	89	GLY	C-N-CA	5.78	131.35	121.87	2	1
1	X	104	PHE	CA-CB-CG	5.75	119.55	113.80	11	1
1	X	74	ARG	NE-CZ-NH2	5.72	124.35	119.20	11	1
1	X	30	ASP	CA-CB-CG	5.60	118.20	112.60	6	2
1	X	42	ILE	CB-CA-C	5.55	119.04	111.88	3	2
1	X	87	ASN	CA-C-N	5.54	127.97	120.38	6	1
1	X	87	ASN	C-N-CA	5.54	127.97	120.38	6	1
1	X	88	ASP	OD1-CG-OD2	-5.51	109.67	122.90	5	1
1	X	63	ASP	CA-CB-CG	5.48	118.08	112.60	8	1
1	X	96	TYR	CA-C-N	5.47	131.99	121.54	8	1
1	X	96	TYR	C-N-CA	5.47	131.99	121.54	8	1
1	X	129	VAL	CA-CB-CG1	5.44	119.65	110.40	8	1
1	X	38	LYS	CA-C-N	5.42	127.80	120.38	5	1
1	X	38	LYS	C-N-CA	5.42	127.80	120.38	5	1
1	X	95	GLU	OE1-CD-OE2	-5.39	109.97	122.90	2	3
1	X	84	ASP	CB-CA-C	5.38	116.84	108.61	11	1
1	X	103	GLN	OE1-CD-NE2	-5.35	117.25	122.60	6	2
1	X	63	ASP	OD1-CG-OD2	-5.35	110.06	122.90	7	2
1	X	67	ASP	OD1-CG-OD2	-5.32	110.14	122.90	8	1
1	X	70	GLU	OE1-CD-OE2	-5.27	110.25	122.90	1	3
1	X	110	ASP	CA-C-N	5.21	131.49	121.54	10	1
1	X	110	ASP	C-N-CA	5.21	131.49	121.54	10	1
1	X	134	HIS	CB-CG-CD2	-5.21	124.43	131.20	11	1
1	X	103	GLN	CA-C-N	5.20	127.50	120.38	11	1
1	X	103	GLN	C-N-CA	5.20	127.50	120.38	11	1
1	X	121	ALA	N-CA-C	5.19	121.86	110.80	3	1
1	X	61	ASP	OD1-CG-OD2	-5.17	110.48	122.90	1	1
1	X	36	PRO	CA-C-N	5.16	127.20	120.28	10	1
1	X	36	PRO	C-N-CA	5.16	127.20	120.28	10	1
1	X	116	ASP	N-CA-C	5.15	115.09	108.34	4	1
1	X	103	GLN	CA-CB-CG	5.14	124.39	114.10	4	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	X	39	ASP	CA-CB-CG	5.13	117.73	112.60	2	1
1	X	66	LEU	CB-CA-C	5.07	119.83	111.31	3	1
1	X	92	ASP	CA-C-N	5.05	127.30	120.38	7	1
1	X	92	ASP	C-N-CA	5.05	127.30	120.38	7	1
1	X	84	ASP	CA-CB-CG	5.01	117.61	112.60	1	1
1	X	131	LEU	CB-CA-C	5.00	119.08	111.02	3	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	X	122	SER	Peptide	9
1	X	41	THR	Peptide	8
1	X	66	LEU	Peptide	6
1	X	120	LEU	Peptide	6
1	X	39	ASP	Peptide	4
1	X	88	ASP	Peptide	4
1	X	40	GLY	Peptide	4
1	X	133	ARG	Peptide,Sidechain	3
1	X	74	ARG	Peptide,Sidechain	3
1	X	30	ASP	Peptide	2
1	X	118	ARG	Sidechain	2
1	X	96	TYR	Sidechain	2
1	X	83	LEU	Peptide	2
1	X	108	ASN	Peptide	2
1	X	35	ASP	Peptide	2
1	X	37	ASP	Peptide	2
1	X	38	LYS	Peptide	2
1	X	121	ALA	Peptide	2
1	X	28	LYS	Peptide	1
1	X	42	ILE	Peptide	1
1	X	58	LEU	Peptide	1
1	X	65	THR	Peptide	1
1	X	110	ASP	Peptide	1
1	X	134	HIS	Peptide	1
1	X	109	PRO	Peptide	1
1	X	130	ASN	Peptide	1
1	X	85	PRO	Peptide	1
1	X	107	ALA	Peptide	1

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	X	796	784	784	2±2
All	All	9588	9408	9408	20

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:X:122:SER:H	1:X:123:PRO:CD	0.60	2.08	4	2
1:X:29:VAL:HG22	1:X:30:ASP:H	0.53	1.62	8	3
1:X:44:LEU:HD23	1:X:104:PHE:CD1	0.50	2.41	6	1
1:X:95:GLU:C	1:X:97:LEU:H	0.46	2.16	8	1
1:X:96:TYR:CZ	1:X:131:LEU:HA	0.46	2.45	8	1
1:X:55:PHE:CD2	1:X:55:PHE:C	0.46	2.94	11	2
1:X:122:SER:H	1:X:123:PRO:HD3	0.45	1.71	4	1
1:X:120:LEU:C	1:X:120:LEU:HD13	0.45	2.36	8	1
1:X:121:ALA:C	1:X:123:PRO:HD2	0.44	2.37	5	1
1:X:128:LEU:O	1:X:128:LEU:HD23	0.44	2.13	1	1
1:X:122:SER:H	1:X:123:PRO:HD2	0.44	1.72	8	1
1:X:57:LYS:HA	1:X:57:LYS:HE2	0.43	1.91	4	1
1:X:41:THR:HG21	1:X:114:THR:OG1	0.42	2.15	5	1
1:X:35:ASP:C	1:X:37:ASP:H	0.42	2.23	8	1
1:X:55:PHE:CZ	1:X:93:LYS:HE3	0.41	2.51	10	1
1:X:44:LEU:C	1:X:44:LEU:HD13	0.40	2.41	2	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	X	107/117 (91%)	86±2 (80±2%)	15±3 (14±3%)	6±2 (6±2%)	2	19
All	All	1284/1404 (91%)	1027 (80%)	180 (14%)	77 (6%)	2	19

All 23 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	X	122	SER	9
1	X	123	PRO	9
1	X	38	LYS	8
1	X	134	HIS	6
1	X	40	GLY	5
1	X	112	ASP	5
1	X	29	VAL	5
1	X	85	PRO	4
1	X	35	ASP	4
1	X	30	ASP	4
1	X	121	ALA	3
1	X	31	ILE	3
1	X	110	ASP	2
1	X	97	LEU	1
1	X	28	LYS	1
1	X	77	GLU	1
1	X	78	ALA	1
1	X	100	VAL	1
1	X	119	GLU	1
1	X	68	ALA	1
1	X	96	TYR	1
1	X	111	ASN	1
1	X	127	ALA	1

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	X	82/92 (89%)	70±3 (85±3%)	12±3 (15±3%)	5	42
All	All	984/1104 (89%)	837 (85%)	147 (15%)	5	42

All 45 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	X	31	ILE	9
1	X	43	ASP	9
1	X	100	VAL	9
1	X	75	VAL	8
1	X	55	PHE	8
1	X	92	ASP	8
1	X	35	ASP	7
1	X	88	ASP	6
1	X	48	LEU	6
1	X	96	TYR	5
1	X	28	LYS	5
1	X	42	ILE	4
1	X	65	THR	4
1	X	70	GLU	4
1	X	97	LEU	4
1	X	105	LYS	4
1	X	120	LEU	4
1	X	118	ARG	4
1	X	30	ASP	3
1	X	131	LEU	3
1	X	58	LEU	3
1	X	74	ARG	3
1	X	128	LEU	3
1	X	103	GLN	2
1	X	86	ASP	2
1	X	38	LYS	1
1	X	45	LYS	1
1	X	116	ASP	1
1	X	110	ASP	1
1	X	119	GLU	1
1	X	122	SER	1
1	X	83	LEU	1
1	X	44	LEU	1
1	X	56	ASP	1
1	X	61	ASP	1
1	X	67	ASP	1
1	X	129	VAL	1
1	X	111	ASN	1
1	X	115	ILE	1
1	X	123	PRO	1
1	X	101	GLU	1

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Mol	Chain	Res	Type	Models (Total)
1	X	108	ASN	1
1	X	29	VAL	1
1	X	84	ASP	1
1	X	126	SER	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 3 ligands modelled in this entry, 3 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

The completeness of assignment taking into account all chemical shift lists is 72% for the well-defined parts and 67% for the entire structure.

7.1 Chemical shift list 1

File name: working_cs.cif

Chemical shift list name: *assigned_chem_shift_list_1*

7.1.1 Bookkeeping

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	998
Number of shifts mapped to atoms	998
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	5

7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	107	2.64 ± 0.18	Should be applied
$^{13}\text{C}_\beta$	100	3.00 ± 0.12	Should be applied
$^{13}\text{C}'$	0	—	None (insufficient data)
^{15}N	100	0.98 ± 0.48	Should be applied

7.1.3 Completeness of resonance assignments

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 72%, i.e. 974 atoms were assigned a chemical shift out of a possible 1356. 0 out of 16 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	401/532 (75%)	203/216 (94%)	102/214 (48%)	96/102 (94%)
Sidechain	566/778 (73%)	380/501 (76%)	186/251 (74%)	0/26 (0%)

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	Total	¹ H	¹³ C	¹⁵ N
Aromatic	7/46 (15%)	6/23 (26%)	1/22 (5%)	0/1 (0%)
Overall	974/1356 (72%)	589/740 (80%)	289/487 (59%)	96/129 (74%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 67%, i.e. 997 atoms were assigned a chemical shift out of a possible 1487. 0 out of 16 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	415/580 (72%)	208/235 (89%)	107/234 (46%)	100/111 (90%)
Sidechain	575/826 (70%)	383/533 (72%)	192/267 (72%)	0/26 (0%)
Aromatic	7/81 (9%)	6/43 (14%)	1/32 (3%)	0/6 (0%)
Overall	997/1487 (67%)	597/811 (74%)	300/533 (56%)	100/143 (70%)

7.1.4 Statistically unusual chemical shifts ⓘ

The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

List Id	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	X	41	THR	HG1	4.78	0.08 – 2.19	17.2
1	X	42	ILE	CG1	14.28	19.24 – 36.26	-7.9
1	X	123	PRO	CD	42.65	45.11 – 55.58	-7.3
1	X	123	PRO	HG3	3.98	0.33 – 3.48	6.6
1	X	93	LYS	HD2	0.35	0.58 – 2.64	-6.1

7.1.5 Random Coil Index (RCI) plots ⓘ

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 X:

