



Full wwPDB NMR Structure Validation Report ⓘ

Mar 25, 2026 – 07:30 AM UTC

PDB ID : 2JP7 / pdb_00002jp7
Title : NMR structure of the Mex67 UBA domain
Authors : Hobeika, M.; Brockmann, C.; Iglesias, N.; Gwizdek, C.; Neuhaus, D.; Stutz, F.; Stewart, M.; Divita, G.; Dargemont, C.
Deposited on : 2007-04-26

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

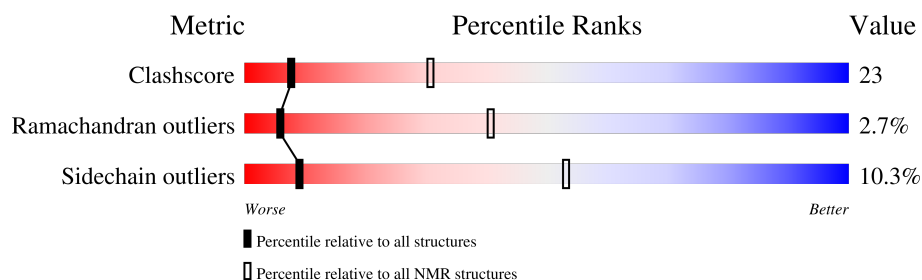
1 Overall quality at a glance

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	57	46% 32% 16% • 5%

2 Ensemble composition and analysis

This entry contains 20 models. Model 5 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *closest to the average*.

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:544-A:597 (54)	0.47	5

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 6 single-model clusters were found.

Cluster number	Models
1	4, 5, 6, 8, 9, 11, 12, 14, 18, 19, 20
2	1, 3, 16
Single-model clusters	2; 7; 10; 13; 15; 17

3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 938 atoms, of which 467 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called mRNA export factor MEX67.

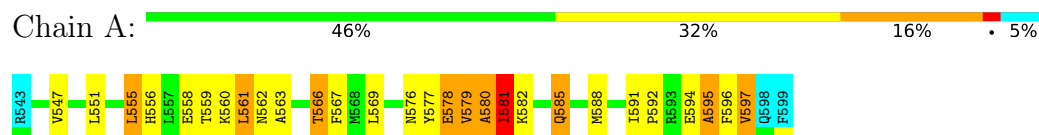
Mol	Chain	Residues	Atoms						Trace
1	A	57	Total	C	H	N	O	S	0
			938	303	467	79	87	2	

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: mRNA export factor MEX67

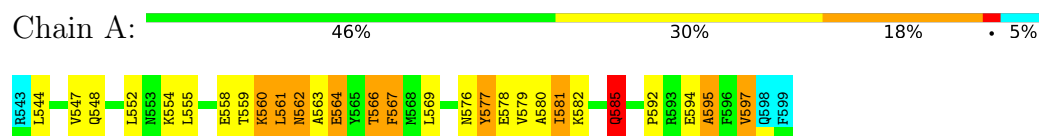


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

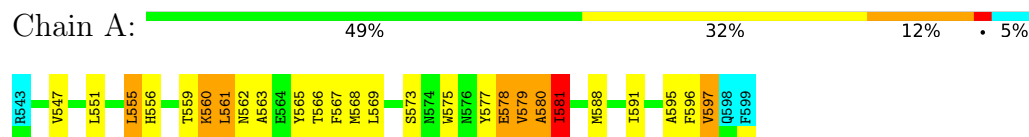
4.2.1 Score per residue for model 1

- Molecule 1: mRNA export factor MEX67



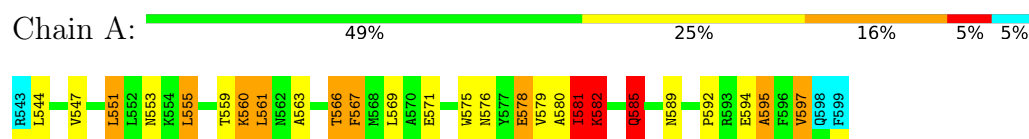
4.2.2 Score per residue for model 2

- Molecule 1: mRNA export factor MEX67



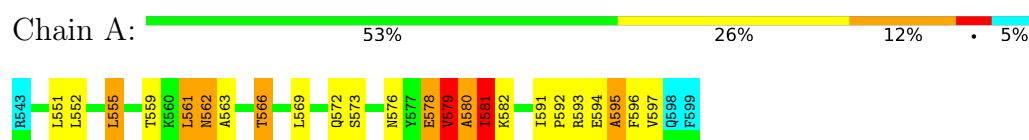
4.2.3 Score per residue for model 3

- Molecule 1: mRNA export factor MEX67



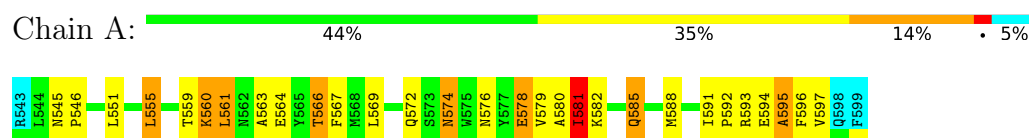
4.2.4 Score per residue for model 4

- Molecule 1: mRNA export factor MEX67



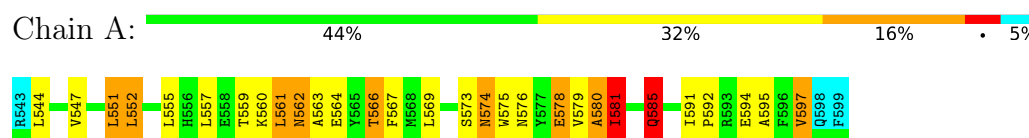
4.2.5 Score per residue for model 5 (medoid)

- Molecule 1: mRNA export factor MEX67



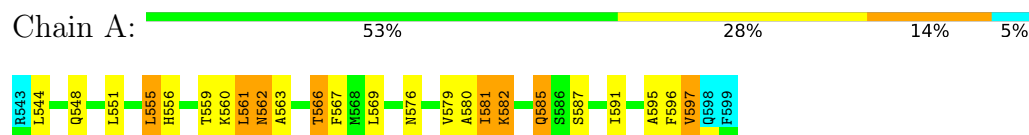
4.2.6 Score per residue for model 6

- Molecule 1: mRNA export factor MEX67



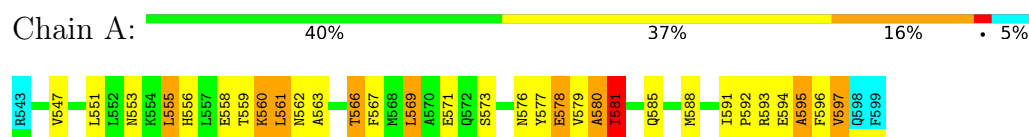
4.2.7 Score per residue for model 7

- Molecule 1: mRNA export factor MEX67



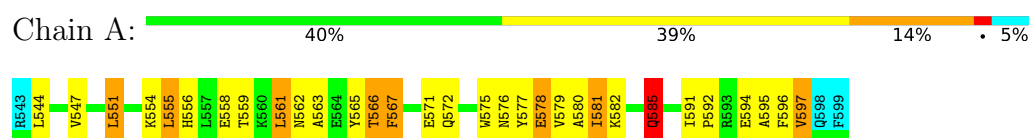
4.2.8 Score per residue for model 8

- Molecule 1: mRNA export factor MEX67



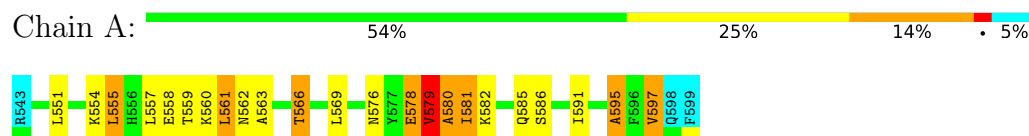
4.2.9 Score per residue for model 9

- Molecule 1: mRNA export factor MEX67



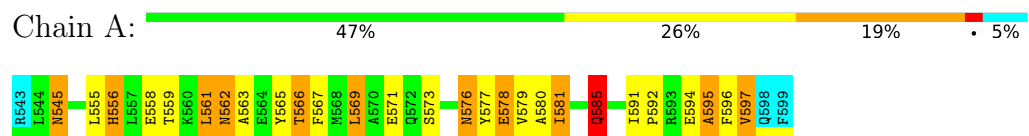
4.2.10 Score per residue for model 10

- Molecule 1: mRNA export factor MEX67



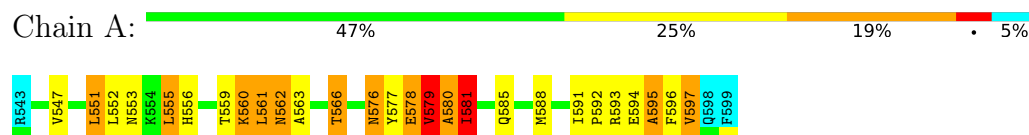
4.2.11 Score per residue for model 11

- Molecule 1: mRNA export factor MEX67



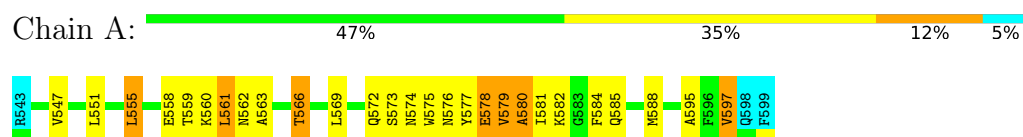
4.2.12 Score per residue for model 12

- Molecule 1: mRNA export factor MEX67



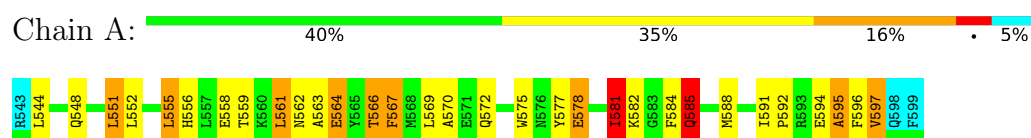
4.2.13 Score per residue for model 13

- Molecule 1: mRNA export factor MEX67



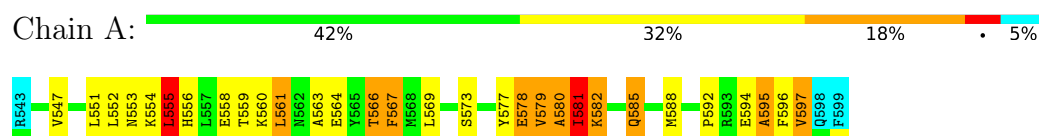
4.2.14 Score per residue for model 14

- Molecule 1: mRNA export factor MEX67



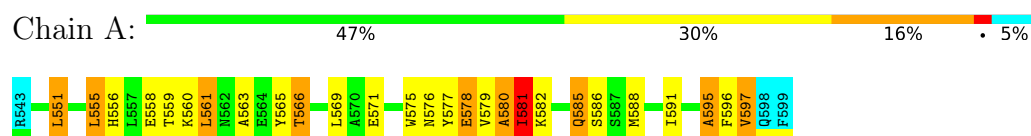
4.2.15 Score per residue for model 15

- Molecule 1: mRNA export factor MEX67



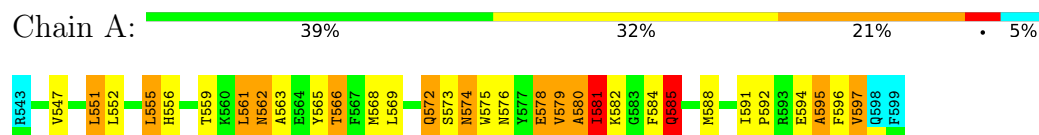
4.2.16 Score per residue for model 16

- Molecule 1: mRNA export factor MEX67



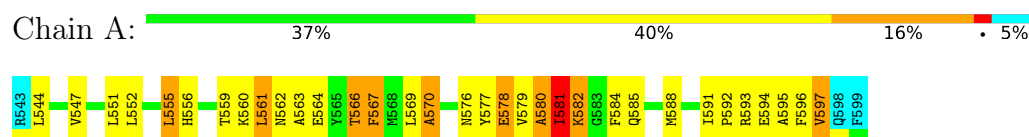
4.2.17 Score per residue for model 17

- Molecule 1: mRNA export factor MEX67



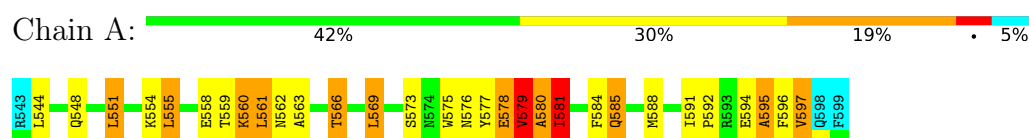
4.2.18 Score per residue for model 18

- Molecule 1: mRNA export factor MEX67



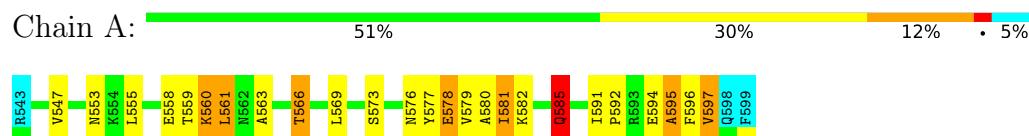
4.2.19 Score per residue for model 19

- Molecule 1: mRNA export factor MEX67



4.2.20 Score per residue for model 20

- Molecule 1: mRNA export factor MEX67



5 Refinement protocol and experimental data overview

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

Of the 200 calculated structures, 20 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
X-PLOR NIH	structure solution	2.13
X-PLOR NIH	refinement	2.13

No chemical shift data was provided.

6 Model quality

6.1 Standard geometry

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	1.75±0.04	10±2/449 (2.2± 0.5%)	1.37±0.04	3±1/609 (0.4± 0.1%)
All	All	1.75	200/8980 (2.2%)	1.37	53/12180 (0.4%)

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	561	LEU	CA-C	-11.94	1.38	1.52	4	20
1	A	581	ILE	N-CA	-9.67	1.34	1.46	2	4
1	A	561	LEU	N-CA	-9.59	1.34	1.46	5	19
1	A	578	GLU	C-N	-8.85	1.23	1.33	12	19
1	A	566	THR	C-O	-8.81	1.14	1.24	19	18
1	A	595	ALA	CA-C	-8.22	1.44	1.53	8	8
1	A	561	LEU	C-N	-8.12	1.21	1.33	10	13
1	A	581	ILE	CA-C	-7.72	1.43	1.52	18	7
1	A	582	LYS	C-O	-7.67	1.15	1.24	17	11
1	A	570	ALA	N-CA	-7.39	1.37	1.46	18	1
1	A	551	LEU	C-O	-7.34	1.15	1.24	6	15
1	A	560	LYS	CA-C	-6.49	1.45	1.53	5	7
1	A	569	LEU	CA-C	-6.48	1.44	1.52	18	1
1	A	567	PHE	C-O	-6.30	1.16	1.24	9	3
1	A	579	VAL	C-N	-6.26	1.24	1.33	13	7
1	A	595	ALA	CA-CB	-6.15	1.44	1.53	15	3
1	A	555	LEU	N-CA	-6.11	1.38	1.46	17	15
1	A	579	VAL	N-CA	-5.98	1.40	1.46	12	1
1	A	554	LYS	C-O	-5.85	1.17	1.24	9	1
1	A	562	ASN	CA-C	-5.78	1.45	1.52	12	5
1	A	556	HIS	ND1-CE1	-5.75	1.26	1.32	14	5
1	A	581	ILE	C-O	-5.69	1.17	1.24	16	1
1	A	595	ALA	C-O	-5.62	1.16	1.24	16	9
1	A	595	ALA	N-CA	-5.17	1.39	1.46	8	2
1	A	581	ILE	CA-CB	5.13	1.60	1.54	17	3
1	A	556	HIS	CG-ND1	-5.06	1.32	1.38	16	2

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	580	ALA	N-CA-C	-7.62	103.44	112.89	2	12
1	A	574	ASN	CA-CB-CG	7.25	119.86	112.60	17	2
1	A	585	GLN	CB-CA-C	6.75	122.17	110.01	9	10
1	A	567	PHE	CA-CB-CG	6.63	120.43	113.80	15	4
1	A	579	VAL	CB-CA-C	6.57	117.09	110.65	12	5
1	A	595	ALA	N-CA-C	-6.06	105.54	113.12	15	3
1	A	566	THR	N-CA-CB	5.54	118.87	110.28	15	4
1	A	585	GLN	CA-CB-CG	5.52	125.15	114.10	1	6
1	A	545	ASN	N-CA-C	-5.46	101.86	108.14	11	1
1	A	561	LEU	N-CA-C	-5.40	101.08	109.72	10	3
1	A	576	ASN	CA-CB-CG	5.13	117.73	112.60	12	1
1	A	591	ILE	N-CA-C	-5.12	104.23	108.63	16	1
1	A	565	TYR	N-CA-C	-5.09	105.43	111.69	16	1

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	439	435	434	20±4
All	All	8780	8700	8680	397

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 23.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:552:LEU:HD13	1:A:567:PHE:HB3	0.79	1.54	15	4
1:A:563:ALA:HA	1:A:566:THR:HG22	0.78	1.54	18	6
1:A:562:ASN:HB2	1:A:595:ALA:HA	0.76	1.55	14	4
1:A:579:VAL:HA	1:A:582:LYS:HG2	0.75	1.57	10	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:573:SER:HB2	1:A:579:VAL:HG13	0.72	1.60	20	7
1:A:592:PRO:HB2	1:A:594:GLU:HG2	0.69	1.64	17	14
1:A:569:LEU:HD13	1:A:580:ALA:HB1	0.69	1.64	17	1
1:A:581:ILE:O	1:A:585:GLN:HB3	0.69	1.88	11	6
1:A:563:ALA:HA	1:A:566:THR:HB	0.67	1.67	17	14
1:A:561:LEU:HA	1:A:595:ALA:O	0.66	1.91	1	19
1:A:581:ILE:O	1:A:585:GLN:HG2	0.66	1.91	10	3
1:A:555:LEU:HD11	1:A:580:ALA:HB3	0.64	1.68	11	2
1:A:555:LEU:O	1:A:559:THR:HG22	0.64	1.92	10	19
1:A:578:GLU:HA	1:A:581:ILE:HB	0.63	1.69	2	18
1:A:581:ILE:O	1:A:585:GLN:HG3	0.63	1.94	6	3
1:A:582:LYS:HA	1:A:585:GLN:NE2	0.63	2.08	7	4
1:A:552:LEU:HD21	1:A:567:PHE:HA	0.62	1.70	6	1
1:A:558:GLU:HG3	1:A:577:TYR:HE1	0.62	1.53	1	1
1:A:562:ASN:CB	1:A:595:ALA:HA	0.62	2.25	9	6
1:A:558:GLU:HG3	1:A:577:TYR:CE1	0.61	2.30	1	2
1:A:581:ILE:O	1:A:585:GLN:HB2	0.61	1.96	16	2
1:A:559:THR:OG1	1:A:561:LEU:HG	0.61	1.96	9	19
1:A:588:MET:HA	1:A:591:ILE:HG12	0.61	1.72	12	5
1:A:578:GLU:C	1:A:581:ILE:HB	0.60	2.22	4	5
1:A:564:GLU:HA	1:A:567:PHE:CE2	0.59	2.32	18	4
1:A:555:LEU:HD21	1:A:569:LEU:HD23	0.58	1.73	2	1
1:A:561:LEU:HD13	1:A:591:ILE:HG21	0.58	1.75	2	5
1:A:582:LYS:O	1:A:586:SER:HB3	0.58	1.98	16	2
1:A:576:ASN:HB3	1:A:579:VAL:CG1	0.57	2.29	1	8
1:A:563:ALA:CA	1:A:566:THR:HG22	0.57	2.29	18	5
1:A:558:GLU:HG3	1:A:577:TYR:OH	0.57	1.99	16	8
1:A:563:ALA:HA	1:A:566:THR:CB	0.57	2.29	2	6
1:A:584:PHE:HE1	1:A:596:PHE:HE1	0.57	1.42	17	1
1:A:588:MET:HB2	1:A:596:PHE:HZ	0.57	1.60	12	4
1:A:544:LEU:HD21	1:A:567:PHE:HE2	0.56	1.61	9	2
1:A:555:LEU:HD21	1:A:580:ALA:HB3	0.55	1.77	10	10
1:A:560:LYS:O	1:A:596:PHE:HA	0.55	2.00	8	2
1:A:555:LEU:HG	1:A:570:ALA:HB2	0.55	1.78	14	1
1:A:561:LEU:O	1:A:565:TYR:HB2	0.54	2.02	2	1
1:A:555:LEU:HD23	1:A:559:THR:HG22	0.54	1.79	11	2
1:A:569:LEU:HA	1:A:572:GLN:HE21	0.54	1.62	5	2
1:A:565:TYR:CD2	1:A:595:ALA:HB2	0.54	2.38	11	3
1:A:567:PHE:O	1:A:571:GLU:HB3	0.53	2.04	3	1
1:A:593:ARG:HA	1:A:596:PHE:HB2	0.53	1.78	5	3
1:A:569:LEU:CD1	1:A:580:ALA:HB1	0.53	2.32	17	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:591:ILE:HB	1:A:596:PHE:CE2	0.52	2.39	11	2
1:A:554:LYS:O	1:A:558:GLU:HG2	0.52	2.04	10	3
1:A:595:ALA:O	1:A:597:VAL:HG13	0.52	2.05	1	11
1:A:556:HIS:HA	1:A:559:THR:OG1	0.52	2.04	2	1
1:A:544:LEU:HB3	1:A:548:GLN:HB2	0.52	1.82	7	4
1:A:551:LEU:HD11	1:A:577:TYR:HB2	0.51	1.83	12	1
1:A:588:MET:HB2	1:A:596:PHE:CZ	0.51	2.41	18	5
1:A:552:LEU:HD21	1:A:567:PHE:CA	0.51	2.34	6	1
1:A:555:LEU:HG	1:A:559:THR:HG23	0.50	1.83	2	1
1:A:573:SER:OG	1:A:580:ALA:HB2	0.50	2.05	2	9
1:A:562:ASN:ND2	1:A:595:ALA:HA	0.50	2.21	1	2
1:A:555:LEU:HD21	1:A:580:ALA:CB	0.50	2.36	10	9
1:A:578:GLU:CA	1:A:581:ILE:HB	0.50	2.35	4	5
1:A:576:ASN:HB2	1:A:579:VAL:HG12	0.50	1.83	12	1
1:A:551:LEU:CD1	1:A:577:TYR:HB2	0.50	2.37	12	1
1:A:569:LEU:HG	1:A:580:ALA:HB1	0.49	1.85	11	8
1:A:563:ALA:O	1:A:566:THR:HB	0.49	2.07	1	4
1:A:561:LEU:HD22	1:A:591:ILE:HG21	0.48	1.85	4	2
1:A:571:GLU:HA	1:A:575:TRP:NE1	0.48	2.22	9	1
1:A:562:ASN:HB2	1:A:597:VAL:CG1	0.48	2.38	19	4
1:A:551:LEU:HD23	1:A:575:TRP:HE3	0.48	1.68	14	9
1:A:565:TYR:O	1:A:568:MET:HG2	0.48	2.09	2	1
1:A:573:SER:HB2	1:A:579:VAL:HG22	0.48	1.85	19	1
1:A:556:HIS:O	1:A:560:LYS:HA	0.48	2.09	15	1
1:A:567:PHE:HB2	1:A:575:TRP:CH2	0.47	2.44	14	1
1:A:594:GLU:HA	1:A:597:VAL:HA	0.47	1.86	15	1
1:A:544:LEU:HD13	1:A:567:PHE:HE1	0.47	1.69	6	1
1:A:591:ILE:HB	1:A:596:PHE:CE1	0.47	2.45	9	2
1:A:556:HIS:HB2	1:A:566:THR:HG21	0.46	1.88	15	4
1:A:561:LEU:CD1	1:A:591:ILE:HG21	0.46	2.40	5	6
1:A:587:SER:O	1:A:591:ILE:HG12	0.45	2.11	7	1
1:A:593:ARG:HA	1:A:596:PHE:HB3	0.45	1.87	4	1
1:A:562:ASN:HB3	1:A:565:TYR:HD2	0.45	1.71	9	1
1:A:588:MET:HA	1:A:591:ILE:CG1	0.45	2.42	12	2
1:A:544:LEU:HD11	1:A:567:PHE:CE2	0.45	2.46	7	1
1:A:581:ILE:HA	1:A:584:PHE:CD2	0.45	2.47	14	1
1:A:552:LEU:HD12	1:A:566:THR:HG23	0.44	1.89	4	3
1:A:576:ASN:HB3	1:A:579:VAL:HG12	0.44	1.88	5	4
1:A:552:LEU:CD1	1:A:567:PHE:HB3	0.44	2.37	18	1
1:A:551:LEU:CD1	1:A:577:TYR:HA	0.44	2.43	2	1
1:A:554:LYS:HE3	1:A:577:TYR:CZ	0.44	2.47	15	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:562:ASN:HB2	1:A:597:VAL:HG11	0.43	1.89	19	2
1:A:555:LEU:HD23	1:A:570:ALA:HB2	0.43	1.90	18	1
1:A:592:PRO:HB2	1:A:594:GLU:CG	0.43	2.44	8	3
1:A:584:PHE:O	1:A:588:MET:HB3	0.43	2.14	18	3
1:A:588:MET:HA	1:A:596:PHE:CE2	0.43	2.48	15	1
1:A:591:ILE:HG22	1:A:595:ALA:HB3	0.43	1.90	10	1
1:A:552:LEU:HD12	1:A:566:THR:HG22	0.43	1.91	17	1
1:A:581:ILE:HD12	1:A:581:ILE:HA	0.42	1.56	6	1
1:A:576:ASN:HB3	1:A:579:VAL:HG11	0.42	1.90	1	1
1:A:567:PHE:O	1:A:571:GLU:HG2	0.42	2.14	8	2
1:A:556:HIS:ND1	1:A:561:LEU:O	0.42	2.51	11	1
1:A:545:ASN:HB2	1:A:546:PRO:HD2	0.42	1.92	5	1
1:A:567:PHE:O	1:A:571:GLU:HB2	0.42	2.14	9	1
1:A:563:ALA:CA	1:A:566:THR:HB	0.42	2.41	17	1
1:A:591:ILE:HB	1:A:596:PHE:HE1	0.42	1.73	9	1
1:A:582:LYS:HA	1:A:585:GLN:HE22	0.41	1.74	7	1
1:A:581:ILE:HA	1:A:581:ILE:HD13	0.41	1.39	12	2
1:A:568:MET:O	1:A:572:GLN:HG3	0.41	2.16	17	1
1:A:551:LEU:HD13	1:A:577:TYR:HA	0.41	1.92	2	1
1:A:582:LYS:NZ	1:A:582:LYS:HB2	0.41	2.29	3	1
1:A:581:ILE:HD13	1:A:581:ILE:HA	0.41	1.35	4	1
1:A:561:LEU:CD2	1:A:591:ILE:HG21	0.41	2.46	10	1
1:A:591:ILE:HB	1:A:596:PHE:HE2	0.41	1.75	7	1
1:A:559:THR:C	1:A:560:LYS:HD2	0.41	2.40	7	3
1:A:563:ALA:HA	1:A:566:THR:OG1	0.41	2.15	2	1
1:A:574:ASN:OD1	1:A:576:ASN:HB2	0.41	2.16	5	1
1:A:569:LEU:CD2	1:A:580:ALA:HB1	0.41	2.46	8	1
1:A:563:ALA:O	1:A:567:PHE:HB2	0.40	2.15	2	1
1:A:566:THR:O	1:A:569:LEU:HB3	0.40	2.16	14	1
1:A:552:LEU:HD23	1:A:575:TRP:CZ3	0.40	2.50	6	1
1:A:561:LEU:HB2	1:A:566:THR:OG1	0.40	2.17	14	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	54/57 (95%)	48±2 (88±3%)	5±2 (9±3%)	1±0 (3±1%)	6	41
All	All	1080/1140 (95%)	953 (88%)	98 (9%)	29 (3%)	6	41

All 7 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	597	VAL	20
1	A	562	ASN	3
1	A	577	TYR	2
1	A	589	ASN	1
1	A	574	ASN	1
1	A	592	PRO	1
1	A	588	MET	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	A	48/51 (94%)	43±2 (90±3%)	5±2 (10±3%)	9	53
All	All	960/1020 (94%)	861 (90%)	99 (10%)	9	53

All 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	581	ILE	19
1	A	585	GLN	13
1	A	547	VAL	12
1	A	569	LEU	8
1	A	576	ASN	8
1	A	579	VAL	6
1	A	560	LYS	5
1	A	553	ASN	5
1	A	564	GLU	4
1	A	572	GLN	4
1	A	574	ASN	3

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Mol	Chain	Res	Type	Models (Total)
1	A	562	ASN	2
1	A	557	LEU	2
1	A	555	LEU	2
1	A	588	MET	1
1	A	582	LYS	1
1	A	552	LEU	1
1	A	593	ARG	1
1	A	545	ASN	1
1	A	571	GLU	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

No chemical shift data were provided