



Full wwPDB EM Validation Report ⓘ

Mar 19, 2026 – 07:41 PM UTC

PDB ID : 9NYY / pdb_00009nyy
EMDB ID : EMD-49946
Title : Nucleic acid bound human SLFN14, State 1
Authors : Van Riper, J.; Martinez-Claros, A.O.; Wang, L.; Schneiderman, H.; Maheshwari, S.; Pillon, M.C.
Deposited on : 2025-03-29
Resolution : 2.73 Å(reported)

This is a Full wwPDB EM 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/EMValidationReportHelp>
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:

EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : **NOT EXECUTED**
MapQ : **FAILED**
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

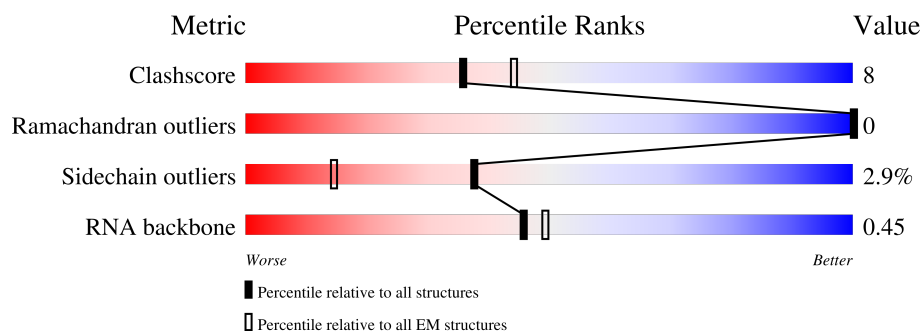
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.73 Å.

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)	EM structures (#Entries)
Clashscore	229148	23984
Ramachandran outliers	224038	23583
Sidechain outliers	223484	23102
RNA backbone	8273	3508

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	943	68% 18% 13%
1	B	943	68% 18% • 13%
2	C	5	20% 80%
3	D	7	57% 43%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 13446 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Protein SLFN14.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	820	Total	C	N	O	S	0	0
			6581	4220	1115	1197	49		
1	B	823	Total	C	N	O	S	0	0
			6607	4236	1119	1203	49		

There are 62 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	913	THR	-	expression tag	UNP P0C7P3
A	914	ARG	-	expression tag	UNP P0C7P3
A	915	THR	-	expression tag	UNP P0C7P3
A	916	ARG	-	expression tag	UNP P0C7P3
A	917	PRO	-	expression tag	UNP P0C7P3
A	918	LEU	-	expression tag	UNP P0C7P3
A	919	GLU	-	expression tag	UNP P0C7P3
A	920	GLN	-	expression tag	UNP P0C7P3
A	921	LYS	-	expression tag	UNP P0C7P3
A	922	LEU	-	expression tag	UNP P0C7P3
A	923	ILE	-	expression tag	UNP P0C7P3
A	924	SER	-	expression tag	UNP P0C7P3
A	925	GLU	-	expression tag	UNP P0C7P3
A	926	GLU	-	expression tag	UNP P0C7P3
A	927	ASP	-	expression tag	UNP P0C7P3
A	928	LEU	-	expression tag	UNP P0C7P3
A	929	ALA	-	expression tag	UNP P0C7P3
A	930	ALA	-	expression tag	UNP P0C7P3
A	931	ASN	-	expression tag	UNP P0C7P3
A	932	ASP	-	expression tag	UNP P0C7P3
A	933	ILE	-	expression tag	UNP P0C7P3
A	934	LEU	-	expression tag	UNP P0C7P3
A	935	ASP	-	expression tag	UNP P0C7P3
A	936	TYR	-	expression tag	UNP P0C7P3
A	937	LYS	-	expression tag	UNP P0C7P3
A	938	ASP	-	expression tag	UNP P0C7P3

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Chain	Residue	Modelled	Actual	Comment	Reference
A	939	ASP	-	expression tag	UNP P0C7P3
A	940	ASP	-	expression tag	UNP P0C7P3
A	941	ASP	-	expression tag	UNP P0C7P3
A	942	LYS	-	expression tag	UNP P0C7P3
A	943	VAL	-	expression tag	UNP P0C7P3
B	913	THR	-	expression tag	UNP P0C7P3
B	914	ARG	-	expression tag	UNP P0C7P3
B	915	THR	-	expression tag	UNP P0C7P3
B	916	ARG	-	expression tag	UNP P0C7P3
B	917	PRO	-	expression tag	UNP P0C7P3
B	918	LEU	-	expression tag	UNP P0C7P3
B	919	GLU	-	expression tag	UNP P0C7P3
B	920	GLN	-	expression tag	UNP P0C7P3
B	921	LYS	-	expression tag	UNP P0C7P3
B	922	LEU	-	expression tag	UNP P0C7P3
B	923	ILE	-	expression tag	UNP P0C7P3
B	924	SER	-	expression tag	UNP P0C7P3
B	925	GLU	-	expression tag	UNP P0C7P3
B	926	GLU	-	expression tag	UNP P0C7P3
B	927	ASP	-	expression tag	UNP P0C7P3
B	928	LEU	-	expression tag	UNP P0C7P3
B	929	ALA	-	expression tag	UNP P0C7P3
B	930	ALA	-	expression tag	UNP P0C7P3
B	931	ASN	-	expression tag	UNP P0C7P3
B	932	ASP	-	expression tag	UNP P0C7P3
B	933	ILE	-	expression tag	UNP P0C7P3
B	934	LEU	-	expression tag	UNP P0C7P3
B	935	ASP	-	expression tag	UNP P0C7P3
B	936	TYR	-	expression tag	UNP P0C7P3
B	937	LYS	-	expression tag	UNP P0C7P3
B	938	ASP	-	expression tag	UNP P0C7P3
B	939	ASP	-	expression tag	UNP P0C7P3
B	940	ASP	-	expression tag	UNP P0C7P3
B	941	ASP	-	expression tag	UNP P0C7P3
B	942	LYS	-	expression tag	UNP P0C7P3
B	943	VAL	-	expression tag	UNP P0C7P3

- Molecule 2 is a RNA chain called RNA (5'-R(P*AP*UP*GP*GP*G)-3').

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	5	Total	C	N	O	P	0	0
			111	49	22	35	5		

- Molecule 3 is a RNA chain called RNA (5'-R(P*CP*CP*CP*AP*CP*UP*C)-3').

Mol	Chain	Residues	Atoms					AltConf	Trace
3	D	7	Total	C	N	O	P	0	0
			142	64	22	49	7		

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
4	A	1	Total	Zn	0
			1	1	
4	B	1	Total	Zn	0
			1	1	

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

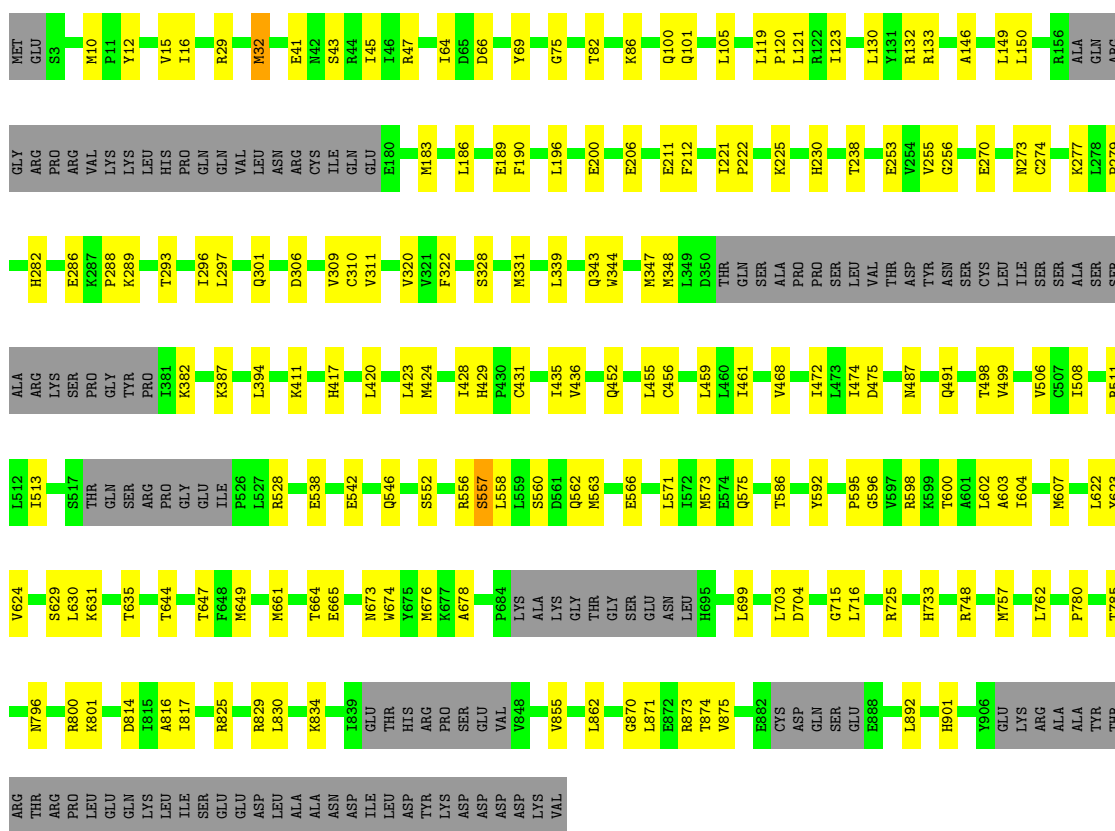
Mol	Chain	Residues	Atoms		AltConf
5	A	2	Total	Mg	0
			2	2	
5	B	1	Total	Mg	0
			1	1	

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-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 outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

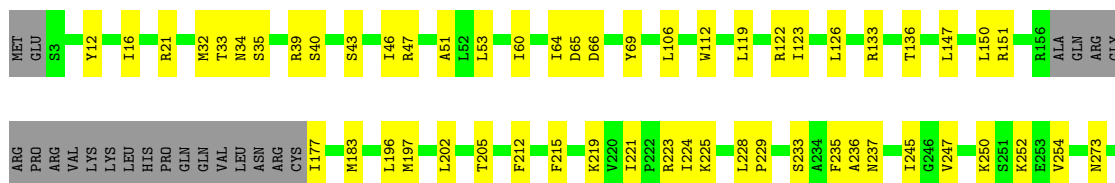
• Molecule 1: Protein SLFN14

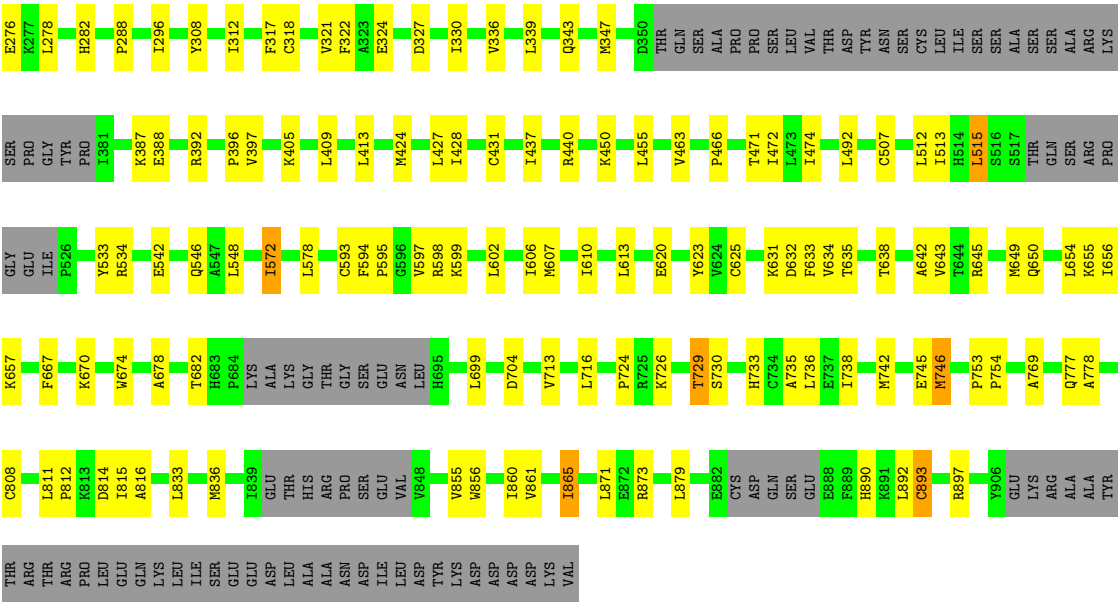
Chain A: 



• Molecule 1: Protein SLFN14

Chain B: 





● Molecule 2: RNA (5'-R(P*AP*UP*GP*GP*G)-3')



● Molecule 3: RNA (5'-R(P*CP*CP*CP*AP*CP*UP*C)-3')



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	338469	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50.2	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOCONTINUUM (6k x 4k)	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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 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	0.13	0/6718	0.32	0/9059
1	B	0.13	0/6744	0.32	0/9094
2	C	0.07	0/124	0.14	0/192
3	D	0.07	0/156	0.14	0/239
All	All	0.13	0/13742	0.31	0/18584

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 the 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 within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	6581	0	6673	99	0
1	B	6607	0	6698	112	0
2	C	111	0	55	2	0
3	D	142	0	77	1	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	2	0	0	0	0
5	B	1	0	0	0	0
All	All	13446	0	13503	210	0

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 8.

All (210) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:282:HIS:HE1	1:B:318:CYS:SG	1.70	1.13
1:A:301:GLN:HG2	1:A:306:ASP:HB2	1.63	0.79
1:A:43:SER:O	1:A:47:ARG:HG3	1.84	0.77
1:A:101:GLN:HE22	1:A:528:ARG:HD3	1.50	0.76
1:B:43:SER:O	1:B:47:ARG:HG3	1.88	0.72
1:B:147:LEU:HA	1:B:150:LEU:HD12	1.72	0.72
1:B:21:ARG:HH12	1:B:65:ASP:HA	1.52	0.72
1:A:387:LYS:HD2	1:A:546:GLN:HB3	1.72	0.71
1:A:221:ILE:HD11	1:A:270:GLU:HG3	1.72	0.70
1:B:730:SER:HB3	1:B:777:GLN:HE22	1.57	0.69
1:A:556:ARG:HH21	1:A:780:PRO:HB3	1.56	0.69
1:B:729:THR:HG21	1:B:769:ALA:HB1	1.75	0.68
1:B:392:ARG:NH1	1:B:392:ARG:HA	2.09	0.68
1:B:215:PHE:HE1	1:B:224:ILE:HD11	1.59	0.68
1:B:595:PRO:HD2	1:B:746:MET:HE1	1.78	0.66
1:B:455:LEU:HD22	1:B:515:LEU:HD21	1.77	0.66
1:A:607:MET:HE1	1:A:623:TYR:HB2	1.76	0.65
1:B:343:GLN:O	1:B:347:MET:HG2	1.97	0.64
1:B:66:ASP:HB3	1:B:69:TYR:HB2	1.79	0.64
1:A:595:PRO:HB3	1:A:892:LEU:HD21	1.79	0.64
1:B:202:LEU:HB2	1:B:252:LYS:HB3	1.80	0.63
1:B:742:MET:O	1:B:746:MET:HG2	1.98	0.62
1:A:343:GLN:O	1:A:347:MET:HG2	2.00	0.62
1:B:147:LEU:O	1:B:151:ARG:HG2	1.99	0.62
1:A:12:TYR:HB2	1:A:16:ILE:HD11	1.82	0.62
1:B:471:THR:HB	1:B:512:LEU:HD23	1.82	0.61
1:B:735:ALA:HB3	1:B:738:ILE:HG13	1.82	0.61
1:B:893:CYS:O	1:B:897:ARG:HG2	2.01	0.61
1:B:424:MET:HE3	1:B:424:MET:HA	1.83	0.61
1:A:382:LYS:HD2	1:A:571:LEU:HD21	1.83	0.60
1:A:575:GLN:HG2	1:A:602:LEU:HD21	1.84	0.60
1:B:21:ARG:NH1	1:B:65:ASP:HA	2.17	0.59
1:A:875:VAL:HG22	1:A:901:HIS:HB2	1.84	0.59
1:A:487:ASN:O	1:A:491:GLN:HG2	2.03	0.59
1:B:392:ARG:HH12	1:B:396:PRO:HA	1.68	0.59
1:A:703:LEU:HD21	1:A:716:LEU:HD11	1.86	0.58
1:B:816:ALA:HB2	1:B:873:ARG:HG3	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:196:LEU:HD12	1:B:197:MET:H	1.69	0.57
1:B:12:TYR:HB2	1:B:16:ILE:HD11	1.85	0.57
1:A:424:MET:HE3	1:A:472:ILE:HD11	1.84	0.57
1:B:598:ARG:HD2	1:B:598:ARG:N	2.20	0.57
1:A:598:ARG:HD3	1:A:733:HIS:HB2	1.87	0.56
1:B:855:VAL:HG21	1:B:871:LEU:HD22	1.88	0.56
1:B:649:MET:O	1:B:650:GLN:HG2	2.06	0.56
1:B:466:PRO:HD3	1:B:534:ARG:HD3	1.87	0.55
1:B:623:TYR:HB3	1:B:642:ALA:HA	1.87	0.55
1:B:254:VAL:HG12	1:B:254:VAL:O	2.07	0.55
1:B:623:TYR:HE2	1:B:625:CYS:HB3	1.71	0.55
1:A:630:LEU:HD22	1:A:870:GLY:HA3	1.89	0.54
1:B:674:TRP:HH2	1:B:716:LEU:HG	1.72	0.54
1:A:814:ASP:OD1	1:A:874:THR:HB	2.07	0.54
1:B:387:LYS:HD2	1:B:546:GLN:HB3	1.89	0.54
1:B:388:GLU:O	1:B:392:ARG:HG2	2.08	0.54
1:A:296:ILE:O	1:A:296:ILE:HD12	2.08	0.54
1:A:596:GLY:HA2	1:A:598:ARG:HH12	1.72	0.54
1:A:796:ASN:HB3	1:A:800:ARG:HH21	1.71	0.54
1:B:427:LEU:HD11	1:B:513:ILE:HG13	1.88	0.54
1:B:33:THR:HG22	1:B:34:ASN:H	1.72	0.54
1:B:229:PRO:HB3	1:B:278:LEU:HD21	1.90	0.53
1:A:631:LYS:O	1:A:635:THR:HG22	2.08	0.53
1:A:123:ILE:HD11	1:A:344:TRP:CD1	2.44	0.52
1:A:297:LEU:HB2	1:A:309:VAL:HG12	1.91	0.52
1:A:498:THR:HG23	1:A:499:VAL:HG23	1.92	0.52
1:A:286:GLU:CD	1:A:286:GLU:H	2.18	0.52
1:B:424:MET:O	1:B:428:ILE:HD12	2.09	0.52
1:A:101:GLN:HA	1:B:177:ILE:HD12	1.91	0.52
1:A:411:LYS:HB3	1:A:411:LYS:NZ	2.24	0.51
1:A:474:ILE:HG13	1:A:475:ASP:N	2.25	0.51
1:B:397:VAL:HG21	1:B:542:GLU:HG3	1.92	0.51
1:B:236:ALA:HB1	1:B:317:PHE:HB2	1.92	0.51
1:B:219:LYS:HB2	1:B:223:ARG:HH21	1.76	0.50
1:A:270:GLU:HA	1:A:273:ASN:ND2	2.27	0.50
1:A:629:SER:HB2	1:A:871:LEU:HD11	1.94	0.50
1:B:60:ILE:O	1:B:106:LEU:HA	2.12	0.50
1:B:607:MET:HE1	1:B:623:TYR:HB2	1.93	0.50
1:A:66:ASP:HB3	1:A:69:TYR:HB2	1.93	0.50
1:B:631:LYS:HD2	1:B:643:VAL:HA	1.94	0.50
1:B:236:ALA:HB3	1:B:321:VAL:HG11	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:440:ARG:HG3	1:B:808:CYS:HA	1.94	0.50
1:A:424:MET:O	1:A:428:ILE:HD12	2.12	0.49
1:B:212:PHE:HB2	1:B:247:VAL:HB	1.95	0.49
1:A:190:PHE:HE1	1:A:196:LEU:HD11	1.77	0.49
1:A:230:HIS:CD2	1:A:230:HIS:H	2.29	0.49
1:B:35:SER:O	1:B:39:ARG:HG2	2.12	0.49
1:A:558:LEU:O	1:A:562:GLN:HG3	2.13	0.49
1:B:745:GLU:HB3	1:B:892:LEU:HD13	1.95	0.49
1:A:816:ALA:HB2	1:A:873:ARG:HD2	1.95	0.48
1:B:196:LEU:HD12	1:B:197:MET:N	2.28	0.48
1:A:221:ILE:HG22	1:A:222:PRO:HD3	1.95	0.48
1:A:557:SER:H	1:A:560:SER:HB2	1.78	0.48
1:A:538:GLU:O	1:A:542:GLU:HG3	2.14	0.48
1:A:801:LYS:HD2	1:A:801:LYS:HA	1.70	0.48
1:A:186:LEU:HA	1:A:189:GLU:HB2	1.95	0.48
1:A:673:ASN:HB3	1:A:676:MET:HB2	1.96	0.48
1:A:678:ALA:HB1	1:A:699:LEU:HD11	1.95	0.48
1:A:394:LEU:HD13	1:A:563:MET:HG2	1.96	0.48
1:B:670:LYS:HA	1:B:670:LYS:HD3	1.63	0.47
1:A:757:MET:HB3	1:A:762:LEU:HD11	1.96	0.47
1:A:423:LEU:HD21	1:A:513:ILE:HG21	1.96	0.47
1:A:603:ALA:O	1:A:607:MET:HG3	2.15	0.47
1:B:606:ILE:O	1:B:610:ILE:HG13	2.15	0.47
1:B:620:GLU:HB3	1:B:657:LYS:HB2	1.96	0.47
1:A:420:LEU:HD11	1:A:472:ILE:HD13	1.97	0.47
1:B:123:ILE:HG21	1:B:347:MET:HE2	1.97	0.47
1:B:133:ARG:HE	1:B:235:PHE:HE1	1.61	0.47
1:A:132:ARG:HB3	1:A:132:ARG:CZ	2.45	0.47
1:B:632:ASP:HA	1:B:635:THR:HG22	1.96	0.47
1:A:15:VAL:HG11	1:A:331:MET:HE2	1.97	0.46
1:A:436:VAL:HA	1:A:459:LEU:O	2.15	0.46
1:B:450:LYS:HE2	1:B:450:LYS:HB2	1.82	0.46
1:B:594:PHE:HD2	1:B:729:THR:HA	1.79	0.46
1:A:279:PRO:HB2	1:A:322:PHE:HB2	1.97	0.46
1:A:596:GLY:HA2	1:A:598:ARG:NH1	2.30	0.46
1:A:674:TRP:NE1	1:A:715:GLY:H	2.13	0.46
1:A:146:ALA:O	1:A:150:LEU:HG	2.15	0.46
1:A:328:SER:HB3	1:A:339:LEU:HD12	1.98	0.46
1:B:196:LEU:HG	1:B:254:VAL:HG11	1.98	0.46
1:B:512:LEU:C	1:B:513:ILE:HD13	2.41	0.46
1:B:593:CYS:HB3	1:B:597:VAL:HG21	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:289:LYS:HE2	1:A:289:LYS:HB3	1.73	0.46
1:A:665:GLU:OE2	1:A:704:ASP:HB3	2.16	0.46
1:B:572:ILE:HG12	1:B:733:HIS:CD2	2.51	0.46
1:A:29:ARG:HA	1:A:32:MET:HE1	1.97	0.45
1:B:46:ILE:HG21	1:B:126:LEU:HD22	1.98	0.45
1:B:392:ARG:HA	1:B:392:ARG:HH11	1.81	0.45
1:B:726:LYS:HB2	1:B:726:LYS:NZ	2.32	0.45
1:A:282:HIS:CE1	1:A:288:PRO:HD2	2.52	0.45
1:B:53:LEU:HD22	1:B:112:TRP:HB2	1.97	0.45
1:B:599:LYS:NZ	1:B:704:ASP:HB2	2.30	0.45
1:B:472:ILE:HD13	1:B:513:ILE:HB	1.98	0.45
1:A:10:MET:HE2	1:A:10:MET:HB3	1.87	0.44
1:A:133:ARG:NH2	1:A:211:GLU:HB2	2.32	0.44
1:A:592:TYR:HE2	1:A:725:ARG:HD2	1.82	0.44
1:B:861:VAL:HG13	1:B:873:ARG:HH21	1.82	0.44
1:B:64:ILE:HG21	1:B:69:TYR:CD2	2.52	0.44
1:B:282:HIS:CE1	1:B:288:PRO:HD2	2.52	0.44
1:A:75:GLY:N	1:B:205:THR:HG21	2.32	0.44
1:B:296:ILE:HG23	1:B:308:TYR:HB3	2.00	0.44
1:B:645:ARG:HD2	1:B:667:PHE:CD1	2.52	0.44
1:A:600:THR:O	1:A:604:ILE:HG12	2.18	0.44
1:B:228:LEU:HD21	1:B:312:ILE:HD13	2.00	0.44
1:B:594:PHE:CD2	1:B:729:THR:HA	2.52	0.44
1:A:468:VAL:HG11	1:A:511:ARG:NH1	2.33	0.44
1:B:833:LEU:HD23	1:B:833:LEU:HA	1.83	0.44
1:A:41:GLU:O	1:A:45:ILE:HG13	2.18	0.44
1:A:435:ILE:HG22	1:A:461:ILE:HB	2.00	0.44
1:A:586:THR:HG22	1:B:724:PRO:HB3	1.99	0.44
1:A:825:ARG:HH12	1:A:829:ARG:HH11	1.66	0.44
1:B:736:LEU:HD23	1:B:778:ALA:HB3	2.00	0.44
1:A:417:HIS:CD2	1:A:455:LEU:HD21	2.53	0.43
1:A:456:CYS:SG	1:A:472:ILE:HD12	2.58	0.43
1:B:512:LEU:O	1:B:513:ILE:HD13	2.18	0.43
1:A:344:TRP:O	1:A:348:MET:HB2	2.18	0.43
1:B:634:VAL:O	1:B:638:THR:HG22	2.18	0.43
1:B:221:ILE:O	1:B:225:LYS:HG3	2.19	0.43
1:A:331:MET:HE3	1:A:331:MET:HB3	1.74	0.43
1:B:330:ILE:HD13	1:B:339:LEU:HD21	2.01	0.43
2:C:4:G:H2'	2:C:5:G:H8	1.84	0.43
1:A:855:VAL:HG21	1:A:871:LEU:HD13	2.00	0.43
1:B:51:ALA:HB2	1:B:336:VAL:HG13	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:817:ILE:O	1:A:862:LEU:HA	2.19	0.43
1:B:678:ALA:HB1	1:B:699:LEU:HD11	2.01	0.43
1:B:865:ILE:HD12	1:B:890:HIS:CE1	2.53	0.43
1:A:674:TRP:HE1	1:A:715:GLY:H	1.67	0.43
1:B:607:MET:CE	1:B:623:TYR:HD1	2.32	0.43
1:B:572:ILE:HG12	1:B:733:HIS:NE2	2.34	0.43
1:B:405:LYS:O	1:B:437:ILE:HA	2.19	0.42
1:A:150:LEU:HD23	1:A:320:VAL:HG21	2.00	0.42
1:A:277:LYS:HA	1:A:277:LYS:HD3	1.69	0.42
2:C:4:G:H2'	2:C:5:G:C8	2.54	0.42
1:B:273:ASN:HA	1:B:276:GLU:HG2	2.00	0.42
1:B:409:LEU:O	1:B:413:LEU:HG	2.20	0.42
1:B:507:CYS:HB3	1:B:533:TYR:HD2	1.84	0.42
1:B:656:ILE:HD13	1:B:656:ILE:HA	1.88	0.42
1:B:812:PRO:HB3	1:B:860:ILE:HG12	2.01	0.42
3:D:3:C:H2'	3:D:4:A:C8	2.54	0.42
1:A:130:LEU:HD11	1:A:149:LEU:HD13	2.02	0.42
1:A:119:LEU:HD12	1:A:120:PRO:HD2	2.02	0.42
1:A:624:VAL:HB	1:A:661:MET:SD	2.60	0.42
1:B:815:ILE:HB	1:B:860:ILE:HG23	2.02	0.42
1:A:64:ILE:HD13	1:A:69:TYR:CD2	2.55	0.42
1:B:250:LYS:HE2	1:B:250:LYS:HA	2.02	0.42
1:B:322:PHE:HB3	1:B:324:GLU:O	2.20	0.42
1:A:834:LYS:HE3	1:A:834:LYS:HB3	1.89	0.41
1:B:753:PRO:HA	1:B:754:PRO:HD3	1.95	0.41
1:B:428:ILE:HD12	1:B:428:ILE:H	1.86	0.41
1:A:119:LEU:HB2	1:B:119:LEU:HD11	2.01	0.41
1:A:200:GLU:N	1:A:200:GLU:OE2	2.53	0.41
1:A:190:PHE:CZ	1:A:196:LEU:HD21	2.56	0.41
1:A:225:LYS:NZ	1:A:274:CYS:HB2	2.35	0.41
1:B:572:ILE:HD13	1:B:572:ILE:HA	1.92	0.41
1:A:573:MET:HE3	1:A:573:MET:HB3	1.95	0.41
1:A:830:LEU:HD23	1:A:830:LEU:HA	1.86	0.41
1:B:233:SER:O	1:B:237:ASN:HB2	2.21	0.41
1:B:474:ILE:H	1:B:474:ILE:HG13	1.67	0.41
1:A:206:GLU:HG2	1:A:212:PHE:CE2	2.56	0.41
1:A:255:VAL:HG23	1:A:256:GLY:H	1.84	0.41
1:B:655:LYS:H	1:B:655:LYS:HD2	1.85	0.41
1:A:644:THR:HG23	1:A:647:THR:H	1.86	0.41
1:B:122:ARG:H	1:B:122:ARG:HG2	1.69	0.41
1:B:126:LEU:HD23	1:B:327:ASP:OD2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:602:LEU:O	1:B:606:ILE:HG12	2.20	0.41
1:B:811:LEU:HD23	1:B:811:LEU:HA	1.87	0.41
1:B:836:MET:HE3	1:B:836:MET:HB3	1.95	0.41
1:A:622:LEU:HD12	1:A:622:LEU:HA	1.94	0.40
1:B:633:PHE:HA	1:B:856:TRP:HE1	1.86	0.40
1:A:82:THR:HG23	1:A:86:LYS:NZ	2.36	0.40
1:A:200:GLU:O	1:A:253:GLU:HA	2.22	0.40
1:A:556:ARG:NH2	1:A:780:PRO:HB3	2.31	0.40
1:B:296:ILE:HD13	1:B:296:ILE:HA	1.89	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 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 EM 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	806/943 (86%)	769 (95%)	37 (5%)	0	100	100
1	B	809/943 (86%)	770 (95%)	39 (5%)	0	100	100
All	All	1615/1886 (86%)	1539 (95%)	76 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 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 EM 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	740/849 (87%)	719 (97%)	21 (3%)	38	60
1	B	743/849 (88%)	721 (97%)	22 (3%)	36	58
All	All	1483/1698 (87%)	1440 (97%)	43 (3%)	38	59

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	MET
1	A	100	GLN
1	A	105	LEU
1	A	121	LEU
1	A	183	MET
1	A	238	THR
1	A	293	THR
1	A	310	CYS
1	A	311	VAL
1	A	429	HIS
1	A	431	CYS
1	A	452	GLN
1	A	506	VAL
1	A	508	ILE
1	A	552	SER
1	A	557	SER
1	A	566	GLU
1	A	649	MET
1	A	664	THR
1	A	748	ARG
1	A	785	THR
1	B	32	MET
1	B	40	SER
1	B	136	THR
1	B	183	MET
1	B	245	ILE
1	B	431	CYS
1	B	463	VAL
1	B	492	LEU
1	B	515	LEU
1	B	548	LEU
1	B	572	ILE
1	B	578	LEU
1	B	613	LEU
1	B	654	LEU

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Mol	Chain	Res	Type
1	B	682	THR
1	B	713	VAL
1	B	729	THR
1	B	746	MET
1	B	814	ASP
1	B	865	ILE
1	B	879	LEU
1	B	893	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	78	GLN
1	A	101	GLN
1	A	298	ASN
1	A	391	GLN
1	A	546	GLN
1	A	577	GLN
1	A	695	HIS
1	B	78	GLN
1	B	291	ASN
1	B	301	GLN
1	B	673	ASN
1	B	803	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	C	4/5 (80%)	2 (50%)	0
3	D	6/7 (85%)	1 (16%)	0
All	All	10/12 (83%)	3 (30%)	0

All (3) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	C	2	U
2	C	3	G
3	D	7	C

There are no RNA pucker outliers to report.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 5 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Map visualisation

This section contains visualisations of the EMDB entry EMD-49946. These allow visual inspection of the internal detail of the map and identification of artifacts.

Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections

This section was not generated.

6.2 Central slices

This section was not generated.

6.3 Largest variance slices

This section was not generated.

6.4 Orthogonal standard-deviation projections (False-color)

This section was not generated.

6.5 Orthogonal surface views

This section was not generated.

6.6 Mask visualisation

This section was not generated. No masks/segmentation were deposited.

7 Map analysis ⓘ

This section contains the results of statistical analysis of the map.

7.1 Map-value distribution ⓘ

This section was not generated.

7.2 Volume estimate versus contour level ⓘ

This section was not generated.

7.3 Rotationally averaged power spectrum ⓘ

This section was not generated. The rotationally averaged power spectrum had issues being displayed.

8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

9 Map-model fit

This section was not generated.