



Full wwPDB EM Validation Report ⓘ

Mar 25, 2026 – 04:30 AM UTC

PDB ID : 8HM0 / pdb_00008hm0
EMDB ID : EMD-34887
Title : F8-A22-E4 complex of MPXV in trimeric form
Authors : Li, Y.N.; Shen, Y.P.; Hu, Z.W.; Yan, R.H.
Deposited on : 2022-12-02
Resolution : 3.10 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

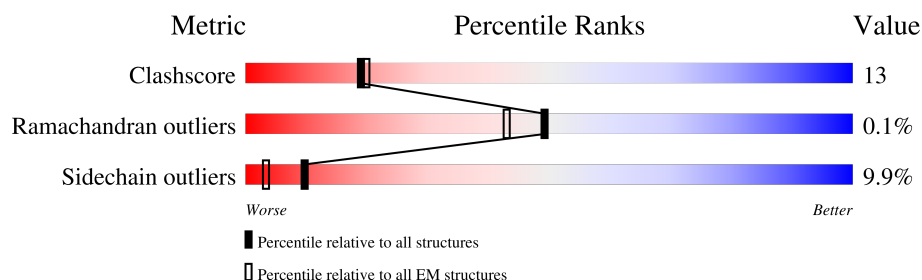
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.10 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

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	C	426	62% 13% • 24%
2	A	1006	61% 29% • 7%
3	B	218	57% 37% 5% •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11331 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 DNA polymerase processivity factor component A20.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	C	323	Total	C	N	O	S	0	0
			1996	1243	361	389	3		

- Molecule 2 is a protein called DNA polymerase.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	A	935	Total	C	N	O	S	0	0
			7592	4845	1270	1425	52		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	108	PHE	LEU	conflict	UNP A0A2L0AR76

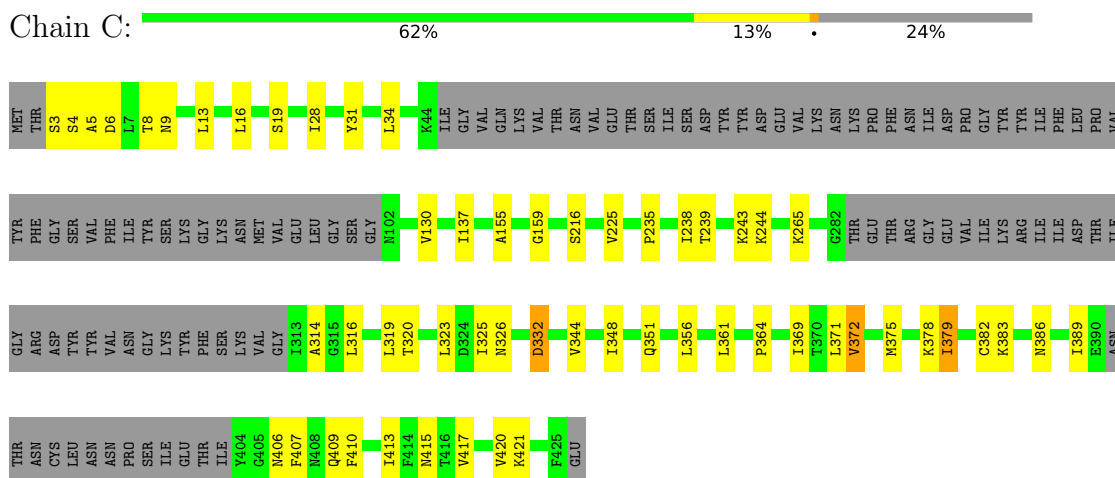
- Molecule 3 is a protein called E4R.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	B	216	Total	C	N	O	S	0	0
			1743	1131	284	323	5		

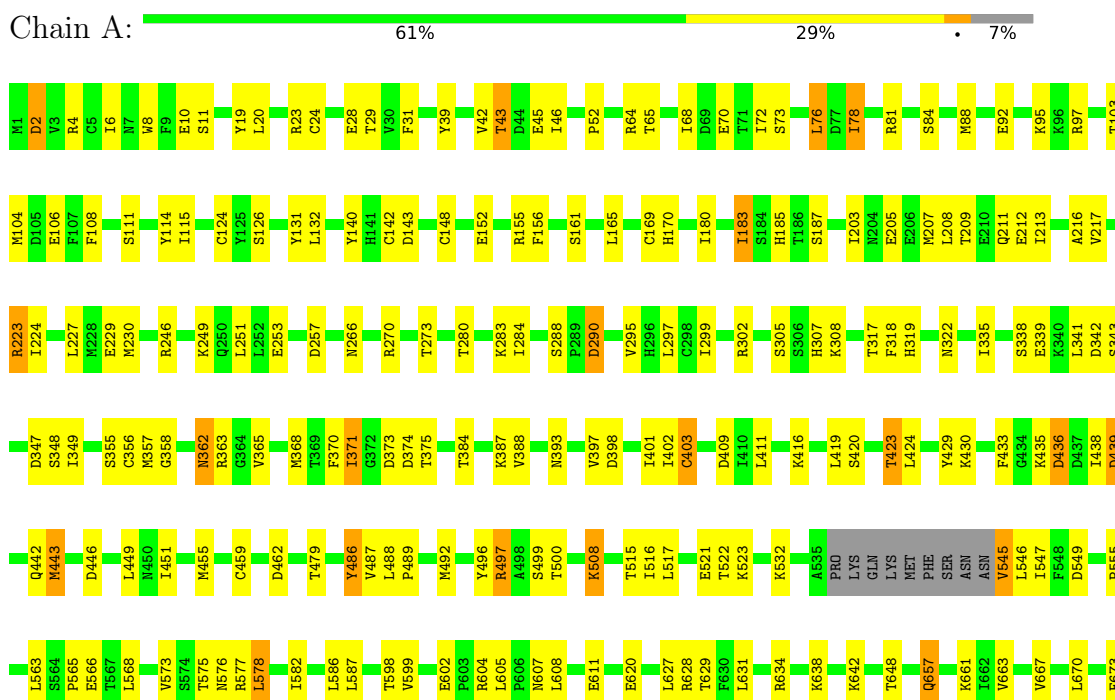
3 Residue-property plots

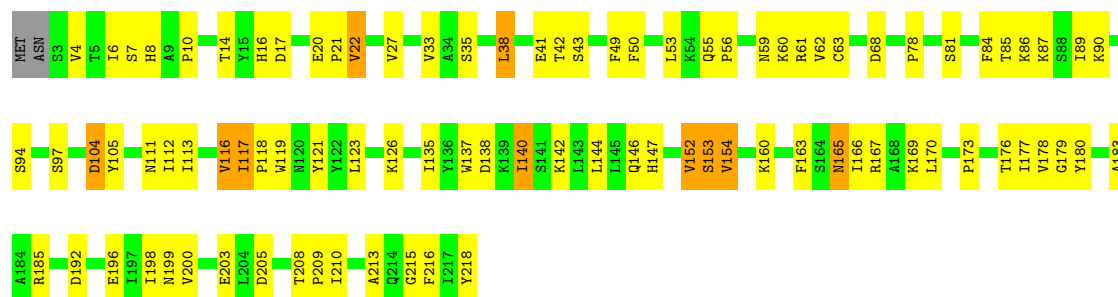
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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: DNA polymerase processivity factor component A20



- Molecule 2: DNA polymerase





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	854700	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1400	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	C	0.16	0/2016	0.44	0/2768
2	A	0.24	0/7745	0.44	0/10451
3	B	0.21	0/1792	0.49	0/2439
All	All	0.22	0/11553	0.45	0/15658

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	C	1996	0	1428	34	0
2	A	7592	0	7560	196	0
3	B	1743	0	1721	55	0
All	All	11331	0	10709	282	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 13.

All (282) 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)
2:A:439:ASP:OD1	2:A:439:ASP:N	2.16	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:573:VAL:HG21	2:A:608:LEU:HD13	1.70	0.74
1:C:371:LEU:HG	1:C:375:MET:HE2	1.70	0.72
2:A:23:ARG:NH1	2:A:29:THR:OG1	2.22	0.72
2:A:8:TRP:HB2	2:A:489:PRO:HB3	1.69	0.72
2:A:187:SER:HG	2:A:459:CYS:HG	1.36	0.72
2:A:439:ASP:HB2	2:A:442:GLN:HB3	1.72	0.72
2:A:738:PRO:HB2	2:A:741:TYR:HE1	1.56	0.70
3:B:8:HIS:HB3	3:B:38:LEU:HD11	1.74	0.70
2:A:205:GLU:OE2	2:A:223:ARG:NH1	2.24	0.69
1:C:9:ASN:HB2	1:C:34:LEU:HD11	1.73	0.69
2:A:81:ARG:HH21	2:A:568:LEU:HB3	1.58	0.68
2:A:115:ILE:HD12	2:A:508:LYS:HG3	1.76	0.67
1:C:225:VAL:HA	1:C:235:PRO:HA	1.77	0.66
2:A:773:GLU:OE2	2:A:776:ARG:NH2	2.27	0.66
2:A:808:THR:O	2:A:826:LYS:NZ	2.27	0.66
2:A:68:ILE:HD13	2:A:517:LEU:HB3	1.78	0.66
2:A:835:VAL:HG12	2:A:836:SER:H	1.61	0.66
2:A:356:CYS:SG	2:A:357:MET:N	2.69	0.65
2:A:709:ASN:HB3	2:A:739:ILE:HA	1.77	0.65
2:A:863:GLN:O	2:A:867:ASP:HB2	1.97	0.65
3:B:60:LYS:HD3	3:B:117:ILE:HB	1.78	0.64
2:A:838:PHE:HB2	2:A:958:PHE:HE2	1.62	0.64
2:A:718:LEU:HG	2:A:782:VAL:HG23	1.80	0.64
2:A:604:ARG:HG3	2:A:605:LEU:HG	1.80	0.64
1:C:239:THR:N	1:C:243:LYS:O	2.27	0.63
2:A:23:ARG:NH2	2:A:257:ASP:OD2	2.31	0.63
2:A:634:ARG:HE	2:A:661:LYS:HB2	1.63	0.63
3:B:105:TYR:HB3	3:B:216:PHE:HD2	1.64	0.63
2:A:305:SER:HB3	2:A:308:LYS:HE3	1.81	0.62
2:A:800:MET:H	2:A:982:LEU:HD21	1.64	0.62
2:A:842:MET:O	2:A:846:TYR:N	2.32	0.62
2:A:989:CYS:SG	2:A:990:ILE:N	2.73	0.61
1:C:378:LYS:HB3	1:C:386:ASN:HB3	1.82	0.61
2:A:602:GLU:O	2:A:685:LYS:NZ	2.28	0.61
3:B:62:VAL:HB	3:B:116:VAL:HG23	1.82	0.61
1:C:238:ILE:HA	1:C:244:LYS:HA	1.82	0.60
3:B:6:ILE:HG22	3:B:27:VAL:HG13	1.83	0.60
2:A:775:GLU:HA	2:A:778:ILE:HG12	1.84	0.60
1:C:361:LEU:HD11	1:C:406:ASN:HB3	1.83	0.59
2:A:810:LYS:HB2	2:A:823:ARG:HG3	1.83	0.59
3:B:87:LYS:HE3	3:B:183:ALA:HA	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:836:SER:O	2:A:890:PHE:HB2	2.02	0.59
2:A:768:ILE:HD13	2:A:791:PHE:HZ	1.67	0.59
2:A:111:SER:O	2:A:115:ILE:HG12	2.03	0.58
1:C:407:PHE:HA	1:C:410:PHE:HB3	1.84	0.58
2:A:185:HIS:HD2	2:A:203:ILE:HG12	1.69	0.58
2:A:52:PRO:O	2:A:97:ARG:NH2	2.30	0.58
2:A:565:PRO:HB3	2:A:680:SER:HB3	1.86	0.58
1:C:372:VAL:HA	1:C:375:MET:HE3	1.85	0.57
2:A:523:LYS:O	2:A:674:ARG:NH2	2.32	0.57
3:B:78:PRO:HB2	3:B:118:PRO:HB2	1.85	0.57
3:B:20:GLU:HG2	3:B:21:PRO:HD3	1.84	0.57
2:A:971:TYR:HB3	2:A:974:ARG:HH21	1.69	0.57
3:B:41:GLU:O	3:B:126:LYS:NZ	2.31	0.57
2:A:563:LEU:HD13	2:A:686:SER:HB3	1.87	0.57
2:A:266:ASN:OD1	2:A:270:ARG:NH2	2.30	0.56
3:B:163:PHE:HA	3:B:166:ILE:HD11	1.87	0.56
2:A:114:TYR:CD2	2:A:487:VAL:HG12	2.40	0.56
2:A:960:LEU:HD13	2:A:966:ILE:HD12	1.87	0.56
3:B:50:PHE:HB3	3:B:53:LEU:HG	1.86	0.56
2:A:397:VAL:HG12	2:A:398:ASP:H	1.70	0.56
2:A:836:SER:OG	2:A:837:LYS:N	2.39	0.56
2:A:720:ASN:HB2	2:A:727:ARG:HH12	1.71	0.55
1:C:375:MET:HG2	1:C:389:ILE:HA	1.87	0.55
2:A:373:ASP:OD1	2:A:375:THR:N	2.37	0.55
2:A:717:THR:OG1	2:A:719:SER:OG	2.24	0.55
2:A:368:MET:HE1	2:A:424:LEU:HB2	1.89	0.55
1:C:344:VAL:O	1:C:348:ILE:HG12	2.06	0.55
2:A:209:THR:HG22	2:A:211:GLN:H	1.71	0.55
2:A:840:LYS:HZ1	2:A:974:ARG:HH12	1.54	0.55
2:A:843:ILE:O	2:A:847:LYS:N	2.37	0.55
1:C:379:ILE:HD12	2:A:582:ILE:HG12	1.89	0.55
2:A:76:LEU:HD11	2:A:587:LEU:HG	1.89	0.55
2:A:227:LEU:HD13	2:A:246:ARG:HB3	1.88	0.55
2:A:841:ASN:ND2	2:A:888:GLU:HB2	2.22	0.54
2:A:973:LYS:O	2:A:977:SER:OG	2.22	0.54
2:A:216:ALA:HB2	2:A:449:LEU:HD11	1.88	0.54
2:A:809:MET:HG3	2:A:810:LYS:H	1.73	0.54
3:B:167:ARG:NH1	3:B:176:THR:OG1	2.32	0.54
2:A:362:ASN:N	2:A:362:ASN:OD1	2.39	0.54
2:A:699:GLU:HG2	2:A:748:VAL:HG22	1.88	0.54
3:B:142:LYS:HG2	3:B:169:LYS:HB2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:883:ARG:HD3	2:A:964:GLN:HE21	1.72	0.54
1:C:417:VAL:HA	1:C:420:VAL:HG12	1.89	0.54
2:A:39:TYR:CE1	2:A:92:GLU:HG2	2.43	0.54
1:C:5:ALA:O	1:C:8:THR:OG1	2.21	0.53
2:A:43:THR:HG22	2:A:88:MET:HE1	1.90	0.53
2:A:187:SER:OG	2:A:459:CYS:SG	2.51	0.53
1:C:13:LEU:HD22	1:C:34:LEU:HB3	1.89	0.53
2:A:891:MET:HE2	2:A:891:MET:HA	1.90	0.53
3:B:4:VAL:HG11	3:B:27:VAL:HG21	1.90	0.53
2:A:545:VAL:HA	2:A:759:ILE:HG22	1.91	0.53
3:B:177:ILE:HD12	3:B:177:ILE:H	1.74	0.52
3:B:22:VAL:HG21	3:B:147:HIS:HB2	1.91	0.52
3:B:154:VAL:HG11	3:B:198:ILE:HG23	1.90	0.52
1:C:369:ILE:HA	1:C:372:VAL:HG12	1.92	0.52
3:B:6:ILE:HD11	3:B:10:PRO:HD2	1.91	0.52
3:B:178:VAL:HG22	3:B:179:GLY:H	1.74	0.52
1:C:369:ILE:HB	2:A:576:ASN:ND2	2.25	0.52
2:A:936:ALA:HB2	2:A:966:ILE:HA	1.92	0.52
1:C:28:ILE:HA	1:C:31:TYR:HB3	1.92	0.51
2:A:577:ARG:HG3	2:A:578:LEU:N	2.25	0.51
2:A:786:ASN:OD1	2:A:786:ASN:N	2.42	0.51
3:B:196:GLU:O	3:B:200:VAL:HG23	2.10	0.51
2:A:782:VAL:HG13	2:A:783:LEU:HD22	1.93	0.51
2:A:70:GLU:OE2	2:A:604:ARG:NH2	2.26	0.51
2:A:843:ILE:HA	2:A:846:TYR:HB3	1.91	0.51
2:A:974:ARG:HD2	2:A:978:GLU:OE2	2.10	0.51
2:A:811:TYR:HE1	2:A:826:LYS:HD3	1.76	0.51
3:B:55:GLN:NE2	3:B:111:ASN:OD1	2.43	0.51
2:A:857:GLY:H	2:A:860:ASN:HB2	1.75	0.51
2:A:620:GLU:OE1	2:A:628:ARG:NH1	2.44	0.51
2:A:446:ASP:OD1	2:A:446:ASP:N	2.40	0.50
3:B:185:ARG:H	3:B:185:ARG:HD2	1.76	0.50
3:B:113:ILE:HB	3:B:116:VAL:HG12	1.93	0.50
2:A:673:PHE:HD2	2:A:676:SER:HB2	1.76	0.50
2:A:735:THR:OG1	2:A:736:SER:N	2.44	0.50
2:A:420:SER:O	2:A:420:SER:OG	2.25	0.50
2:A:825:ASN:CG	2:A:827:GLY:H	2.20	0.50
2:A:717:THR:HA	2:A:731:PRO:HG3	1.93	0.49
2:A:714:PHE:CZ	2:A:778:ILE:HG22	2.48	0.49
2:A:104:MET:HG3	2:A:516:ILE:HG23	1.94	0.49
2:A:604:ARG:N	2:A:611:GLU:OE1	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:23:ARG:HH22	2:A:257:ASP:CG	2.20	0.49
2:A:850:LEU:HA	2:A:853:MET:HG2	1.94	0.49
2:A:24:CYS:HB2	2:A:28:GLU:HG2	1.95	0.49
1:C:216:SER:O	1:C:265:LYS:HA	2.12	0.49
3:B:16:HIS:CE1	3:B:56:PRO:HB3	2.47	0.49
2:A:403:CYS:HB2	2:A:419:LEU:HG	1.95	0.49
1:C:130:VAL:CB	1:C:137:ILE:O	2.60	0.49
1:C:409:GLN:O	1:C:413:ILE:HG22	2.13	0.48
2:A:497:ARG:H	2:A:497:ARG:HE	1.61	0.48
2:A:838:PHE:HZ	2:A:883:ARG:HE	1.59	0.48
2:A:373:ASP:OD1	2:A:375:THR:OG1	2.28	0.48
2:A:573:VAL:HG23	2:A:611:GLU:H	1.79	0.48
3:B:167:ARG:HH22	3:B:173:PRO:HA	1.78	0.48
2:A:733:VAL:HG12	2:A:734:LYS:H	1.78	0.48
2:A:741:TYR:HE2	2:A:766:LYS:HD2	1.79	0.48
2:A:880:PHE:HA	2:A:964:GLN:HB2	1.96	0.48
2:A:853:MET:SD	2:A:864:VAL:HG23	2.53	0.48
2:A:721:PRO:HD2	2:A:784:PHE:CZ	2.49	0.48
3:B:146:GLN:HG2	3:B:170:LEU:HD23	1.96	0.48
2:A:856:GLU:H	2:A:860:ASN:HD22	1.61	0.48
2:A:284:ILE:HB	2:A:297:LEU:HB2	1.95	0.47
2:A:840:LYS:HZ2	2:A:974:ARG:HH22	1.62	0.47
2:A:682:ALA:O	2:A:686:SER:HB2	2.14	0.47
2:A:2:ASP:OD1	2:A:2:ASP:N	2.46	0.47
2:A:124:CYS:SG	2:A:155:ARG:HA	2.54	0.47
3:B:7:SER:OG	3:B:8:HIS:N	2.47	0.47
1:C:325:ILE:HG13	1:C:326:ASN:N	2.30	0.47
2:A:549:ASP:OD1	2:A:790:GLU:HB2	2.14	0.47
2:A:555:PRO:HA	2:A:627:LEU:HD13	1.96	0.47
3:B:61:ARG:HH21	3:B:210:ILE:HD13	1.79	0.47
2:A:577:ARG:HG3	2:A:578:LEU:H	1.80	0.47
3:B:56:PRO:HG2	3:B:59:ASN:HD21	1.78	0.47
2:A:290:ASP:OD1	2:A:290:ASP:N	2.38	0.47
2:A:11:SER:HB2	2:A:19:TYR:HE2	1.79	0.47
2:A:767:SER:HA	2:A:770:ILE:HG22	1.96	0.47
1:C:319:LEU:HD22	1:C:413:ILE:HD11	1.96	0.47
2:A:6:ILE:HD11	2:A:23:ARG:HB2	1.96	0.47
2:A:384:THR:O	2:A:388:VAL:HG23	2.15	0.47
3:B:42:THR:HG21	3:B:49:PHE:HZ	1.80	0.47
2:A:299:ILE:CG2	2:A:318:PHE:HB3	2.45	0.46
2:A:143:ASP:OD1	2:A:143:ASP:N	2.36	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:363:ARG:HG2	2:A:368:MET:HE3	1.96	0.46
2:A:371:ILE:HD12	2:A:416:LYS:HB2	1.96	0.46
2:A:563:LEU:HD12	2:A:687:CYS:HB3	1.98	0.46
2:A:866:ILE:HG21	2:A:870:ARG:HH21	1.80	0.46
2:A:486:TYR:HE1	2:A:663:VAL:HB	1.80	0.46
3:B:42:THR:OG1	3:B:43:SER:N	2.46	0.46
1:C:332:ASP:OD1	1:C:332:ASP:N	2.47	0.46
2:A:705:ALA:HA	2:A:744:ARG:HA	1.98	0.46
2:A:398:ASP:N	2:A:398:ASP:OD1	2.47	0.46
2:A:72:ILE:HD12	2:A:604:ARG:HD2	1.97	0.46
3:B:123:LEU:HD11	3:B:144:LEU:HD12	1.98	0.46
2:A:45:GLU:HG2	2:A:46:ILE:N	2.30	0.46
2:A:78:ILE:H	2:A:78:ILE:HG12	1.57	0.45
2:A:185:HIS:CD2	2:A:203:ILE:HG12	2.51	0.45
1:C:16:LEU:HD13	1:C:31:TYR:HB2	1.97	0.45
2:A:398:ASP:HA	2:A:430:LYS:HE3	1.97	0.45
3:B:213:ALA:HA	3:B:216:PHE:CE1	2.52	0.45
2:A:19:TYR:OH	2:A:290:ASP:OD2	2.26	0.45
2:A:825:ASN:OD1	2:A:827:GLY:N	2.45	0.45
2:A:890:PHE:HE2	2:A:892:LEU:HD23	1.81	0.45
3:B:55:GLN:HE22	3:B:111:ASN:HD21	1.65	0.45
3:B:154:VAL:HG13	3:B:177:ILE:HD13	1.97	0.45
2:A:960:LEU:HD11	2:A:968:TYR:OH	2.17	0.45
2:A:627:LEU:O	2:A:631:LEU:HG	2.17	0.45
3:B:203:GLU:OE1	3:B:209:PRO:HG3	2.17	0.45
2:A:169:CYS:SG	2:A:183:ILE:HG22	2.57	0.45
2:A:224:ILE:HG22	2:A:229:GLU:HB3	1.99	0.45
2:A:317:THR:OG1	2:A:319:HIS:NE2	2.42	0.45
3:B:63:CYS:SG	3:B:152:VAL:HG21	2.57	0.45
3:B:119:TRP:HZ2	3:B:144:LEU:HB3	1.83	0.44
2:A:305:SER:OG	2:A:307:HIS:O	2.33	0.44
2:A:709:ASN:OD1	2:A:735:THR:HG23	2.16	0.44
2:A:853:MET:HA	2:A:863:GLN:NE2	2.32	0.44
2:A:992:PHE:HA	2:A:995:ARG:HG2	1.98	0.44
2:A:170:HIS:HB2	2:A:207:MET:HE1	2.00	0.44
3:B:205:ASP:OD1	3:B:205:ASP:N	2.49	0.44
2:A:4:ARG:NH1	2:A:156:PHE:O	2.45	0.44
2:A:370:PHE:HE2	2:A:429:TYR:HB2	1.83	0.44
2:A:20:LEU:O	2:A:31:PHE:HA	2.18	0.44
2:A:545:VAL:HG12	2:A:758:GLU:HA	2.00	0.44
2:A:973:LYS:HA	2:A:976:THR:HG22	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:990:ILE:O	2:A:994:GLN:HG2	2.18	0.44
3:B:104:ASP:OD1	3:B:104:ASP:N	2.32	0.44
3:B:198:ILE:HD12	3:B:198:ILE:H	1.83	0.44
3:B:216:PHE:HB2	3:B:218:TYR:CZ	2.53	0.44
2:A:766:LYS:HA	2:A:766:LYS:HD3	1.72	0.43
3:B:81:SER:OG	3:B:89:ILE:HG21	2.19	0.43
1:C:3:SER:OG	1:C:6:ASP:OD2	2.32	0.43
2:A:809:MET:HG3	2:A:810:LYS:N	2.33	0.43
2:A:521:GLU:OE2	2:A:522:THR:N	2.46	0.43
2:A:103:THR:OG1	2:A:104:MET:N	2.51	0.43
2:A:810:LYS:HD3	2:A:810:LYS:HA	1.75	0.43
2:A:423:THR:O	2:A:424:LEU:HD23	2.19	0.43
2:A:436:ASP:C	2:A:438:ILE:HD12	2.43	0.43
2:A:573:VAL:HG22	2:A:608:LEU:HD22	2.00	0.43
3:B:153:SER:O	3:B:154:VAL:HG23	2.18	0.43
2:A:358:GLY:HA2	2:A:371:ILE:O	2.18	0.43
1:C:4:SER:O	1:C:8:THR:HG23	2.18	0.43
1:C:383:LYS:HB3	1:C:383:LYS:HE2	1.79	0.43
2:A:106:GLU:H	2:A:106:GLU:HG2	1.66	0.43
2:A:401:ILE:HG22	2:A:403:CYS:SG	2.59	0.43
2:A:638:LYS:HD3	2:A:657:GLN:HG3	2.01	0.43
3:B:167:ARG:NH2	3:B:173:PRO:HA	2.33	0.43
2:A:642:LYS:HB2	2:A:642:LYS:HE3	1.80	0.42
2:A:64:ARG:HE	2:A:515:THR:HG22	1.83	0.42
2:A:108:PHE:CE2	2:A:500:THR:HG23	2.54	0.42
2:A:575:THR:HA	2:A:608:LEU:HD23	2.01	0.42
2:A:867:ASP:O	2:A:871:SER:OG	2.25	0.42
3:B:137:TRP:O	3:B:138:ASP:C	2.62	0.42
2:A:842:MET:HE1	2:A:972:PHE:HE1	1.83	0.42
2:A:64:ARG:NE	2:A:515:THR:HG22	2.35	0.42
2:A:357:MET:HE3	2:A:357:MET:HB3	1.88	0.42
3:B:113:ILE:HG12	3:B:215:GLY:HA3	2.00	0.42
1:C:314:ALA:HB3	1:C:409:GLN:HB3	2.00	0.42
2:A:811:TYR:CE1	2:A:826:LYS:HD3	2.55	0.42
2:A:844:LYS:HA	2:A:847:LYS:HE2	2.01	0.42
2:A:451:ILE:HG22	2:A:455:MET:HE3	2.02	0.42
2:A:685:LYS:HE3	2:A:685:LYS:HB3	1.84	0.42
2:A:339:GLU:HG3	2:A:341:LEU:HD13	2.01	0.42
2:A:862:ASN:HA	2:A:865:CYS:SG	2.60	0.42
2:A:751:ASP:OD1	2:A:751:ASP:N	2.53	0.42
3:B:165:ASN:OD1	3:B:165:ASN:N	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:335:ILE:HG21	2:A:349:ILE:HD13	2.02	0.42
2:A:828:THR:HG22	2:A:828:THR:O	2.20	0.42
2:A:114:TYR:HD2	2:A:487:VAL:HG12	1.84	0.41
2:A:840:LYS:NZ	2:A:974:ARG:HH12	2.16	0.41
2:A:213:ILE:O	2:A:217:VAL:HG23	2.20	0.41
2:A:443:MET:HB3	2:A:455:MET:HE2	2.02	0.41
3:B:140:ILE:H	3:B:140:ILE:HG12	1.45	0.41
1:C:351:GLN:O	1:C:421:LYS:NZ	2.43	0.41
2:A:488:LEU:HD23	2:A:488:LEU:HA	1.94	0.41
3:B:86:LYS:O	3:B:90:LYS:HE2	2.21	0.41
1:C:19:SER:O	1:C:19:SER:OG	2.32	0.41
1:C:3:SER:HA	3:B:192:ASP:HA	2.03	0.41
1:C:155:ALA:O	1:C:159:GLY:N	2.54	0.41
2:A:132:LEU:HB3	2:A:140:TYR:HB3	2.03	0.41
2:A:373:ASP:CG	2:A:374:ASP:N	2.79	0.41
2:A:607:ASN:OD1	2:A:607:ASN:N	2.52	0.41
2:A:749:TYR:CE2	2:A:756:PHE:HB2	2.56	0.41
1:C:383:LYS:HE3	1:C:415:ASN:HB3	2.03	0.41
2:A:131:TYR:HD2	2:A:148:CYS:HB2	1.85	0.41
2:A:532:LYS:HA	2:A:532:LYS:HD2	1.84	0.41
2:A:809:MET:HG3	2:A:823:ARG:HA	2.03	0.41
3:B:94:SER:O	3:B:97:SER:OG	2.31	0.41
3:B:178:VAL:HG22	3:B:179:GLY:N	2.35	0.41
2:A:132:LEU:HD23	2:A:142:CYS:HB2	2.03	0.40
2:A:283:LYS:HE3	2:A:283:LYS:HB2	1.79	0.40
2:A:342:ASP:OD1	2:A:343:SER:N	2.53	0.40
2:A:695:ILE:O	2:A:699:GLU:HG3	2.20	0.40
2:A:249:LYS:O	2:A:253:GLU:HG3	2.21	0.40
2:A:492:MET:HE3	2:A:496:TYR:HE2	1.87	0.40
2:A:706:GLU:OE1	2:A:770:ILE:HD11	2.21	0.40
2:A:778:ILE:O	2:A:783:LEU:HD23	2.22	0.40
3:B:121:TYR:OH	3:B:169:LYS:HD3	2.22	0.40
2:A:251:LEU:HD23	2:A:251:LEU:HA	1.95	0.40
2:A:860:ASN:O	2:A:864:VAL:HG12	2.21	0.40
3:B:196:GLU:HA	3:B:199:ASN:ND2	2.36	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 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 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	C	315/426 (74%)	290 (92%)	24 (8%)	1 (0%)	36	67
2	A	927/1006 (92%)	830 (90%)	97 (10%)	0	100	100
3	B	214/218 (98%)	192 (90%)	22 (10%)	0	100	100
All	All	1456/1650 (88%)	1312 (90%)	143 (10%)	1 (0%)	49	78

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	364	PRO

5.3.2 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 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	C	118/396 (30%)	110 (93%)	8 (7%)	14	42
2	A	856/928 (92%)	770 (90%)	86 (10%)	7	29
3	B	195/200 (98%)	173 (89%)	22 (11%)	5	23
All	All	1169/1524 (77%)	1053 (90%)	116 (10%)	10	29

All (116) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	316	LEU
1	C	320	THR

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Mol	Chain	Res	Type
1	C	323	LEU
1	C	332	ASP
1	C	356	LEU
1	C	372	VAL
1	C	379	ILE
1	C	382	CYS
2	A	2	ASP
2	A	10	GLU
2	A	42	VAL
2	A	43	THR
2	A	65	THR
2	A	73	SER
2	A	76	LEU
2	A	78	ILE
2	A	84	SER
2	A	95	LYS
2	A	126	SER
2	A	152	GLU
2	A	161	SER
2	A	165	LEU
2	A	180	ILE
2	A	183	ILE
2	A	208	LEU
2	A	212	GLU
2	A	223	ARG
2	A	230	MET
2	A	273	THR
2	A	280	THR
2	A	288	SER
2	A	290	ASP
2	A	295	VAL
2	A	302	ARG
2	A	322	ASN
2	A	338	SER
2	A	347	ASP
2	A	348	SER
2	A	355	SER
2	A	362	ASN
2	A	365	VAL
2	A	371	ILE
2	A	387	LYS
2	A	393	ASN

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Mol	Chain	Res	Type
2	A	402	ILE
2	A	403	CYS
2	A	409	ASP
2	A	411	LEU
2	A	423	THR
2	A	433	PHE
2	A	435	LYS
2	A	436	ASP
2	A	439	ASP
2	A	443	MET
2	A	462	ASP
2	A	479	THR
2	A	486	TYR
2	A	497	ARG
2	A	499	SER
2	A	508	LYS
2	A	545	VAL
2	A	546	LEU
2	A	547	ILE
2	A	566	GLU
2	A	578	LEU
2	A	586	LEU
2	A	598	THR
2	A	599	VAL
2	A	629	THR
2	A	648	THR
2	A	657	GLN
2	A	667	VAL
2	A	670	LEU
2	A	686	SER
2	A	709	ASN
2	A	717	THR
2	A	735	THR
2	A	748	VAL
2	A	749	TYR
2	A	752	THR
2	A	757	THR
2	A	775	GLU
2	A	783	LEU
2	A	784	PHE
2	A	820	VAL
2	A	830	GLU

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Mol	Chain	Res	Type
2	A	836	SER
2	A	839	HIS
2	A	845	THR
2	A	887	LEU
2	A	977	SER
2	A	980	VAL
2	A	983	LEU
2	A	988	LEU
3	B	14	THR
3	B	17	ASP
3	B	22	VAL
3	B	33	VAL
3	B	35	SER
3	B	38	LEU
3	B	68	ASP
3	B	84	PHE
3	B	85	THR
3	B	104	ASP
3	B	112	ILE
3	B	116	VAL
3	B	117	ILE
3	B	135	ILE
3	B	140	ILE
3	B	152	VAL
3	B	153	SER
3	B	154	VAL
3	B	160	LYS
3	B	165	ASN
3	B	180	TYR
3	B	208	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	136	ASN
2	A	185	HIS
2	A	265	HIS
2	A	322	ASN
2	A	336	GLN
2	A	442	GLN
2	A	450	ASN
2	A	562	ASN

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Mol	Chain	Res	Type
2	A	801	GLN
2	A	863	GLN
3	B	25	GLN
3	B	55	GLN
3	B	59	ASN
3	B	111	ASN
3	B	146	GLN
3	B	188	GLN
3	B	206	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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](#)

There are no ligands in this entry.

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-34887. 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.