



Full wwPDB EM Validation Report ⓘ

Mar 20, 2026 – 12:27 AM UTC

PDB ID : 8YEO / pdb_00008yeo
EMDB ID : EMD-39200
Title : Type I-FHNH Cascade-dsDNA R-loop complex
Authors : Li, Z.
Deposited on : 2024-02-22
Resolution : 3.44 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

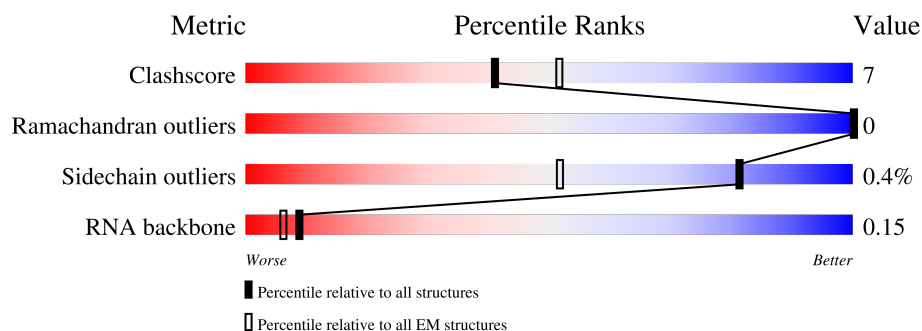
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.44 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.



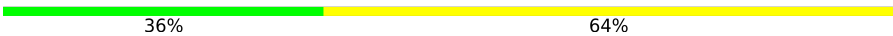
Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	229148	23984
Ramachandran outliers	224038	23583
Sidechain outliers	223484	23102
RNA backbone	8273	3508

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	255	78% (green), 22% (yellow)
2	J	344	70% (green), 28% (yellow), .. (red/orange)
3	B	181	69% (green), 28% (yellow), .. (red/orange)
4	D	335	73% (green), 15% (yellow), 12% (grey)
4	E	335	84% (green), 13% (yellow), . (grey)
4	F	335	85% (green), 11% (yellow), . (grey)
4	G	335	79% (green), 18% (yellow), . (grey)

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Mol	Chain	Length	Quality of chain
4	H	335	 83% 13% .
4	I	335	 86% 10% .
5	C	60	 35% 53% 12%
6	T	48	 71% 29%
7	N	11	 36% 64%

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 24197 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 Cas5f.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	255	Total	C	N	O	S	0	0
			2010	1281	338	377	14		

- Molecule 2 is a protein called Cas8f fusion with HNH.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	J	340	Total	C	N	O	S	0	0
			2718	1719	468	518	13		

- Molecule 3 is a protein called Cas6f.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	B	175	Total	C	N	O	S	0	0
			1415	913	239	258	5		

- Molecule 4 is a protein called Cas7f.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	I	324	Total	C	N	O	S	0	0
			2639	1688	436	500	15		
4	D	295	Total	C	N	O	S	0	0
			2412	1546	397	456	13		
4	E	322	Total	C	N	O	S	0	0
			2621	1677	433	496	15		
4	F	323	Total	C	N	O	S	0	0
			2630	1682	434	499	15		
4	G	323	Total	C	N	O	S	0	0
			2630	1682	434	499	15		
4	H	323	Total	C	N	O	S	0	0
			2630	1682	434	499	15		

- Molecule 5 is a RNA chain called 60-nt crRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	C	60	Total	C	N	O	P	0	0
			1287	577	240	411	59		

- Molecule 6 is DNA/RNA hybrid called TS.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	T	48	Total	C	N	O	P	0	0
			975	468	168	291	48		

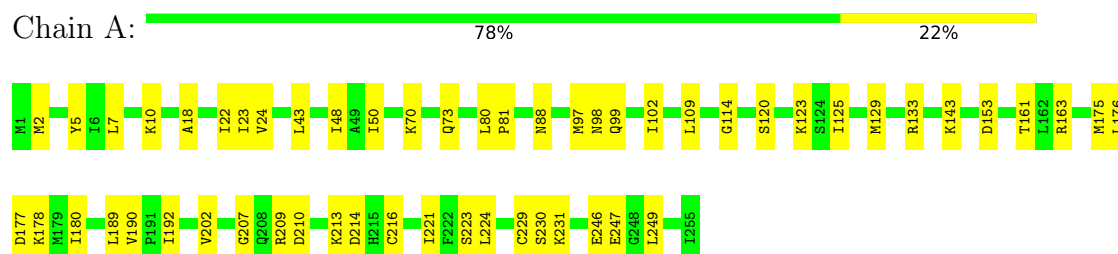
- Molecule 7 is DNA/RNA hybrid called NTS.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	N	11	Total	C	N	O	P	0	0
			230	109	44	66	11		

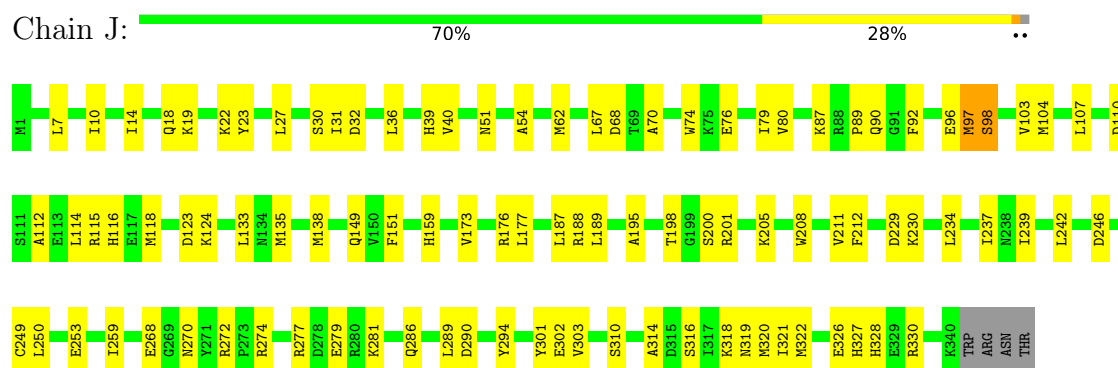
3 Residue-property plots

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.

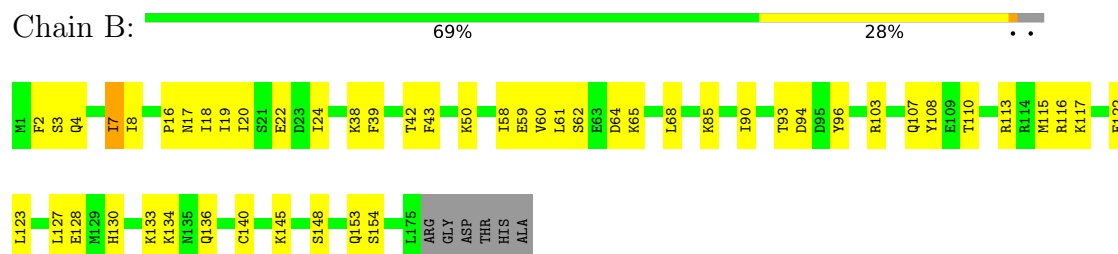
- Molecule 1: Cas5f



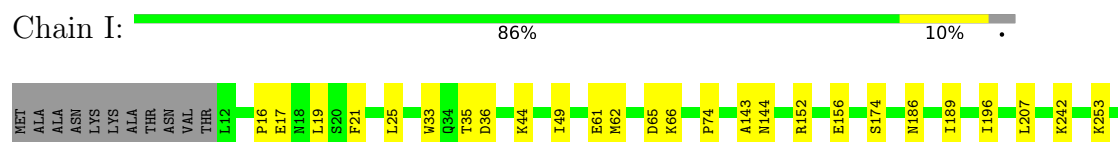
- Molecule 2: Cas8f fusion with HNH



- Molecule 3: Cas6f



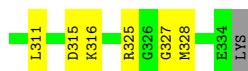
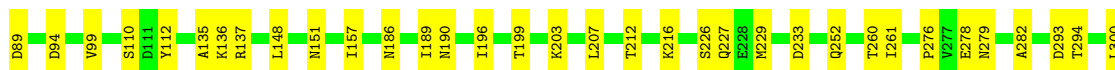
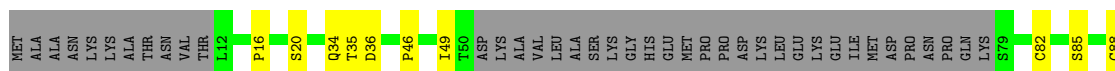
- Molecule 4: Cas7f





- Molecule 4: Cas7f

Chain D: 73% 15% 12%



- Molecule 4: Cas7f

Chain E: 84% 13% .



- Molecule 4: Cas7f

Chain F: 85% 11% .



- Molecule 4: Cas7f

Chain G: 79% 18% .



- Molecule 4: Cas7f

Y206	Y215	M229	D232	D235	K236	M249	K253	R259	T260	I261	Y265	F273	P274	I275	M279	F290	R303	E309	K331	E334	LYS															
ALA	ASN	LYS	LYS	THR	VAL	THR	L12	K13	S14	R15	F21	K58	I71	M72	D73	P74	N75	P76	Q77	D89	K103	D111	Y114	D123	K136	R137	N140	N144	R152	I158	D166	S174	F177	N186	I189	F205

U1	U2	U3	U4	U5	U6	U7	U8	U9	U10	U11	U12	U13	U14	U15	U16	U17	U18	U19	U20	U21	U22	U23	U24	U25	U26	U27	U28	U29	U30	U31	U32	U33	U34	U35	U36	U37	U38	U39	U40	U41	U42	U43	U44	U45	U46	U47	U48	U49	U50	U51	U52	U53	U54	U55	U56	U57	U58	U59	U60	U61	U62	U63	U64	U65	U66	U67	U68	U69	U70	U71	U72	U73	U74	U75	U76	U77	U78	U79	U80	U81	U82	U83	U84	U85	U86	U87	U88	U89	U90	U91	U92	U93	U94	U95	U96	U97	U98	U99	U100
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Node Type	Number of Nodes
G6	100
C7	100
C8	100
A9	100
A10	100
G11	100
T17	100
C20	100
C26	100
C27	100
T28	100
T32	100
A33	100
A38	100
G46	100
C47	100
T48	100
A49	100
C53	100

G9	T12
	T13
	A14
	G15
	C16
	G17
	G18
	A19

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	48756	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	54	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	2200	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	A	0.12	0/2045	0.37	0/2756
2	J	0.14	0/2773	0.38	0/3745
3	B	0.17	0/1446	0.43	0/1945
4	D	0.12	0/2465	0.33	0/3327
4	E	0.11	0/2680	0.31	0/3618
4	F	0.12	0/2689	0.32	0/3630
4	G	0.11	0/2689	0.31	0/3630
4	H	0.10	0/2689	0.27	0/3630
4	I	0.10	0/2698	0.27	0/3641
5	C	0.15	0/1443	0.35	0/2250
6	T	0.20	0/1090	0.45	0/1678
7	N	0.15	0/258	0.34	0/397
All	All	0.13	0/24965	0.34	0/34247

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	2010	0	2030	41	0
2	J	2718	0	2708	72	0
3	B	1415	0	1437	35	0
4	D	2412	0	2379	31	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	E	2621	0	2596	27	0
4	F	2630	0	2602	28	0
4	G	2630	0	2602	42	0
4	H	2630	0	2602	29	0
4	I	2639	0	2615	23	0
5	C	1287	0	645	24	0
6	T	975	0	545	16	0
7	N	230	0	125	4	0
All	All	24197	0	22886	326	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 7.

All (326) 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)
6:T:8:DC:H2''	6:T:9:DA:H5'	1.64	0.80
2:J:103:VAL:O	2:J:107:LEU:HB2	1.93	0.69
4:I:74:PRO:HG2	6:T:38:A:H1'	1.76	0.68
1:A:50:ILE:HG22	1:A:114:GLY:HA3	1.77	0.66
1:A:80:LEU:HD12	1:A:81:PRO:HD2	1.78	0.66
2:J:286:GLN:NE2	2:J:290:ASP:OD2	2.27	0.65
4:H:71:ILE:HG23	4:H:72:MET:HG2	1.80	0.64
1:A:43:LEU:HD22	1:A:48:ILE:HG13	1.80	0.64
2:J:79:ILE:HG22	2:J:90:GLN:HE21	1.64	0.63
4:G:106:MET:HE1	4:G:115:ARG:HG3	1.78	0.63
2:J:274:ARG:HG2	2:J:277:ARG:HH21	1.62	0.63
1:A:229:CYS:SG	1:A:230:SER:N	2.72	0.63
3:B:103:ARG:NH2	5:C:60:G:N7	2.47	0.63
6:T:28:DT:H5'	4:G:77:GLN:HE22	1.64	0.63
4:F:156:GLU:HG2	4:F:157:ILE:HG13	1.80	0.62
4:H:186:ASN:HB3	4:H:189:ILE:HB	1.81	0.62
2:J:40:VAL:HG21	2:J:80:VAL:HG21	1.82	0.62
1:A:98:ASN:HA	2:J:189:LEU:HD11	1.82	0.62
2:J:205:LYS:HZ3	7:N:19:A:HO3'	1.45	0.62
3:B:42:THR:HB	3:B:59:GLU:HB3	1.82	0.62
1:A:177:ASP:OD1	2:J:176:ARG:NH1	2.33	0.62
6:T:20:C:H1'	4:F:74:PRO:HG2	1.82	0.61
4:I:49:ILE:HD12	4:I:242:LYS:HG2	1.82	0.61
4:D:136:LYS:NZ	4:D:190:ASN:OD1	2.33	0.60
4:F:58:LYS:HD2	4:F:60:HIS:HE1	1.67	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:112:TYR:HB2	4:E:288:ILE:HD11	1.85	0.59
2:J:289:LEU:HB3	2:J:294:TYR:HB2	1.84	0.59
2:J:314:ALA:O	2:J:319:ASN:ND2	2.36	0.58
2:J:112:ALA:O	2:J:116:HIS:ND1	2.32	0.58
4:D:229:MET:HG2	4:E:53:ALA:HB1	1.85	0.58
4:E:15:ARG:O	4:E:333:GLN:NE2	2.36	0.58
1:A:88:ASN:ND2	6:T:46:G:OP2	2.37	0.58
3:B:108:TYR:HE1	3:B:128:GLU:HG3	1.69	0.58
4:G:16:PRO:HG2	4:G:19:LEU:HB2	1.86	0.57
3:B:133:LYS:O	3:B:136:GLN:HB3	2.04	0.56
1:A:175:MET:HB3	2:J:67:LEU:HD13	1.86	0.56
4:I:311:LEU:O	4:I:316:LYS:NZ	2.38	0.56
3:B:110:THR:HA	3:B:113:ARG:HH12	1.68	0.56
5:C:51:G:N2	5:C:55:A:OP2	2.39	0.56
2:J:253:GLU:OE2	2:J:327:HIS:NE2	2.35	0.56
3:B:50:LYS:NZ	3:B:140:CYS:SG	2.79	0.56
3:B:130:HIS:O	3:B:133:LYS:HB2	2.06	0.56
4:E:65:ASP:O	4:E:69:LYS:NZ	2.38	0.56
4:E:159:GLU:HG2	4:E:214:PHE:HB2	1.88	0.56
4:D:135:ALA:HB1	4:D:196:ILE:HG23	1.89	0.55
4:D:279:ASN:OD1	4:D:325:ARG:NH2	2.40	0.55
2:J:149:GLN:OE1	5:C:4:A:N6	2.40	0.55
4:G:71:ILE:HG23	4:G:72:MET:HG2	1.88	0.55
4:F:288:ILE:HD12	4:F:289:PRO:HD2	1.89	0.55
4:H:76:PRO:O	4:H:77:GLN:NE2	2.39	0.55
2:J:274:ARG:HA	2:J:277:ARG:HB2	1.88	0.55
5:C:11:G:O2'	4:H:21:PHE:O	2.25	0.55
2:J:27:LEU:HD13	2:J:89:PRO:HG3	1.88	0.55
4:G:156:GLU:HG2	4:G:157:ILE:HG13	1.89	0.55
2:J:89:PRO:HA	2:J:92:PHE:HB3	1.89	0.54
4:E:262:ASP:HB3	4:E:275:ILE:HG13	1.89	0.54
1:A:97:MET:O	1:A:99:GLN:NE2	2.40	0.54
4:E:156:GLU:HG2	4:E:157:ILE:HG13	1.87	0.54
4:H:89:ASP:OD1	4:H:89:ASP:N	2.36	0.54
2:J:7:LEU:O	2:J:19:LYS:NZ	2.40	0.54
1:A:209:ARG:NH2	2:J:149:GLN:O	2.41	0.54
4:I:65:ASP:OD1	4:I:65:ASP:N	2.39	0.54
4:H:75:ASN:OD1	4:H:77:GLN:NE2	2.41	0.54
1:A:73:GLN:OE1	1:A:73:GLN:N	2.39	0.54
4:E:146:ARG:NH2	4:E:179:LEU:O	2.39	0.54
4:F:65:ASP:OD1	4:F:65:ASP:N	2.38	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:14:ILE:HG13	2:J:19:LYS:HG3	1.89	0.54
4:I:253:LYS:NZ	5:C:10:A:OP1	2.40	0.53
4:E:115:ARG:NH1	4:E:119:TYR:OH	2.41	0.53
4:G:262:ASP:HB3	4:G:275:ILE:HG13	1.91	0.53
4:D:260:THR:HA	4:D:276:PRO:HA	1.90	0.53
1:A:161:THR:OG1	1:A:163:ARG:NH1	2.42	0.53
4:G:136:LYS:NZ	4:G:190:ASN:OD1	2.42	0.53
4:E:304:MET:HB2	4:E:311:LEU:HD11	1.91	0.53
5:C:22:A:O2'	5:C:23:U:O4'	2.26	0.53
4:F:144:ASN:O	4:F:152:ARG:NH1	2.42	0.53
4:H:232:ASP:O	4:H:236:LYS:NZ	2.42	0.53
3:B:7:ILE:HG23	3:B:85:LYS:HB3	1.91	0.53
2:J:110:ASP:OD1	2:J:115:ARG:NH2	2.42	0.53
3:B:148:SER:HB2	3:B:153:GLN:H	1.75	0.52
3:B:18:ILE:O	3:B:22:GLU:HB2	2.09	0.52
4:F:15:ARG:NH2	4:G:61:GLU:OE1	2.41	0.52
3:B:148:SER:N	3:B:153:GLN:O	2.38	0.52
5:C:14:G:OP2	4:H:259:ARG:NH2	2.41	0.52
4:G:16:PRO:HB2	4:G:110:SER:HB3	1.92	0.52
4:H:136:LYS:O	4:H:140:ASN:ND2	2.42	0.52
1:A:178:LYS:HE2	2:J:62:MET:HE1	1.92	0.52
4:H:144:ASN:O	4:H:152:ARG:NH1	2.42	0.52
2:J:279:GLU:O	2:J:281:LYS:NZ	2.43	0.51
1:A:120:SER:HA	1:A:123:LYS:HD2	1.93	0.51
4:H:158:ILE:HG23	4:H:215:VAL:HG22	1.93	0.51
1:A:210:ASP:OD2	1:A:213:LYS:NZ	2.42	0.51
4:F:120:GLN:NE2	4:F:124:GLU:OE2	2.43	0.51
2:J:76:GLU:OE2	2:J:201:ARG:NH1	2.43	0.51
2:J:195:ALA:O	2:J:198:THR:OG1	2.28	0.51
4:D:252:GLN:HE21	4:E:54:VAL:HG12	1.76	0.51
4:G:94:ASP:OD1	4:G:94:ASP:N	2.43	0.51
4:G:242:LYS:NZ	4:G:245:GLY:O	2.44	0.51
4:G:260:THR:HA	4:G:276:PRO:HA	1.92	0.51
4:H:303:ARG:NH2	4:H:309:GLU:OE2	2.42	0.51
2:J:277:ARG:HD3	2:J:310:SER:HA	1.93	0.50
4:G:156:GLU:O	4:G:175:LYS:NZ	2.44	0.50
3:B:39:PHE:HA	3:B:61:LEU:O	2.12	0.50
1:A:190:VAL:HG23	1:A:192:ILE:HD11	1.93	0.50
4:D:311:LEU:O	4:D:316:LYS:NZ	2.44	0.50
4:F:143:ALA:O	4:F:174:SER:OG	2.27	0.50
1:A:176:LEU:HD21	2:J:176:ARG:HG3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:233:ASP:OD1	4:D:233:ASP:N	2.37	0.50
4:I:21:PHE:O	5:C:5:G:O2'	2.27	0.50
1:A:246:GLU:HB2	1:A:249:LEU:HB2	1.94	0.49
4:G:144:ASN:O	4:G:152:ARG:NH1	2.45	0.49
4:G:136:LYS:HD3	4:G:184:GLU:HG3	1.94	0.49
5:C:10:A:O2'	5:C:11:G:O4'	2.28	0.49
5:C:14:G:N3	4:H:253:LYS:NZ	2.60	0.49
2:J:200:SER:OG	2:J:201:ARG:N	2.45	0.49
3:B:65:LYS:HE3	3:B:90:ILE:HB	1.94	0.49
3:B:116:ARG:NH1	3:B:122:GLU:OE1	2.45	0.49
4:F:16:PRO:HG2	4:F:19:LEU:HB2	1.94	0.49
4:G:40:ARG:NE	4:G:159:GLU:OE2	2.45	0.49
4:G:105:SER:OG	4:G:105:SER:O	2.30	0.49
4:I:186:ASN:HB3	4:I:189:ILE:HB	1.95	0.49
5:C:37:A:H1'	4:E:56:ALA:HB2	1.94	0.49
4:D:16:PRO:HB2	4:D:110:SER:HB3	1.95	0.49
4:I:16:PRO:HG2	4:I:19:LEU:HB2	1.93	0.49
4:E:31:LYS:NZ	4:E:244:GLU:OE1	2.45	0.49
1:A:2:MET:SD	1:A:2:MET:N	2.85	0.49
4:E:294:THR:HA	4:E:299:LYS:HG3	1.95	0.49
7:N:16:C:H2''	7:N:17:G:C8	2.47	0.49
4:D:49:ILE:HD13	4:D:82:CYS:HB2	1.95	0.48
4:D:89:ASP:OD1	4:D:89:ASP:N	2.46	0.48
4:D:94:ASP:OD1	4:D:212:THR:OG1	2.26	0.48
1:A:202:VAL:HG12	1:A:216:CYS:HB2	1.95	0.48
4:H:137:ARG:NH1	4:H:261:ILE:O	2.44	0.48
2:J:322:MET:SD	2:J:322:MET:N	2.86	0.48
4:G:33:TRP:HB2	4:G:92:ARG:HB3	1.96	0.48
4:F:156:GLU:O	4:F:175:LYS:NZ	2.47	0.48
1:A:133:ARG:O	5:C:2:U:N3	2.41	0.48
2:J:302:GLU:HB2	2:J:328:HIS:HD2	1.79	0.48
4:H:174:SER:HA	4:H:177:PHE:HD2	1.79	0.48
2:J:39:HIS:HB2	2:J:54:ALA:HB3	1.96	0.48
2:J:104:MET:SD	2:J:104:MET:N	2.86	0.48
2:J:32:ASP:N	2:J:32:ASP:OD1	2.38	0.48
3:B:8:ILE:HG21	3:B:16:PRO:HB3	1.96	0.48
4:F:35:THR:OG1	4:F:36:ASP:N	2.47	0.48
4:D:226:SER:OG	4:D:227:GLN:N	2.47	0.48
4:G:35:THR:O	4:G:90:THR:OG1	2.31	0.48
4:D:293:ASP:OD1	4:D:294:THR:N	2.47	0.47
4:D:282:ALA:HB1	4:E:72:MET:HE1	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:E:136:LYS:O	4:E:140:ASN:ND2	2.46	0.47
1:A:10:LYS:HB2	1:A:143:LYS:HB2	1.97	0.47
1:A:223:SER:OG	1:A:224:LEU:N	2.47	0.47
2:J:90:GLN:HG2	2:J:201:ARG:HB3	1.96	0.47
2:J:70:ALA:H	2:J:74:TRP:HE1	1.63	0.47
6:T:8:DC:H2''	6:T:9:DA:H2'	1.97	0.47
4:F:311:LEU:O	4:F:316:LYS:NZ	2.47	0.47
4:H:235:ASP:OD1	4:H:236:LYS:N	2.47	0.47
4:I:144:ASN:O	4:I:152:ARG:NH1	2.47	0.47
4:F:252:GLN:HE21	4:G:54:VAL:HG12	1.79	0.47
2:J:51:ASN:OD1	2:J:51:ASN:N	2.48	0.47
4:G:49:ILE:HD12	4:G:242:LYS:HG2	1.97	0.47
4:G:58:LYS:HE3	4:G:73:ASP:HB2	1.96	0.47
4:D:203:LYS:HA	4:D:203:LYS:HD3	1.70	0.47
4:G:99:VAL:HG23	4:G:199:THR:HG21	1.97	0.46
2:J:230:LYS:HD3	2:J:230:LYS:HA	1.55	0.46
3:B:107:GLN:NE2	5:C:46:C:OP1	2.48	0.46
5:C:35:U:H1'	4:D:20:SER:HB3	1.98	0.46
4:G:139:VAL:HA	4:G:142:ILE:HG22	1.96	0.46
4:G:173:ASN:OD1	4:G:176:SER:OG	2.30	0.46
4:E:49:ILE:HD12	4:E:242:LYS:HG2	1.98	0.46
4:F:265:TYR:OH	4:F:273:PHE:O	2.33	0.46
2:J:246:ASP:HB2	2:J:249:CYS:HB3	1.97	0.46
4:D:278:GLU:OE1	4:E:57:SER:OG	2.34	0.46
4:G:234:ASP:OD1	4:G:234:ASP:N	2.49	0.46
2:J:250:LEU:O	2:J:301:TYR:OH	2.34	0.46
5:C:43:G:H1'	5:C:44:U:C2	2.51	0.46
4:G:73:ASP:OD1	4:G:73:ASP:N	2.49	0.46
4:H:111:ASP:OD2	4:H:114:TYR:N	2.47	0.46
7:N:14:A:H2''	7:N:15:G:C8	2.51	0.46
1:A:231:LYS:HB3	1:A:231:LYS:HE2	1.80	0.45
2:J:27:LEU:O	2:J:31:ILE:HG12	2.15	0.45
4:I:62:MET:HG2	4:I:66:LYS:HE2	1.98	0.45
6:T:8:DC:C2'	6:T:9:DA:H2'	2.46	0.45
4:I:33:TRP:CD2	4:I:44:LYS:HD2	2.51	0.45
4:F:293:ASP:OD1	4:F:293:ASP:N	2.47	0.45
2:J:36:LEU:HD22	2:J:135:MET:HG2	1.98	0.45
4:D:85:SER:OG	4:D:88:CYS:SG	2.69	0.45
1:A:70:LYS:HD2	1:A:70:LYS:HA	1.84	0.45
4:F:332:LYS:HE3	4:F:334:GLU:HG2	1.98	0.45
1:A:177:ASP:HA	1:A:180:ILE:HG12	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:62:SER:OG	3:B:64:ASP:O	2.34	0.45
3:B:134:LYS:HB3	3:B:134:LYS:HE2	1.79	0.45
4:I:156:GLU:OE1	4:H:206:TYR:OH	2.28	0.45
4:D:186:ASN:HB3	4:D:189:ILE:HB	1.98	0.45
5:C:39:A:H1'	5:C:40:A:C8	2.52	0.45
4:F:49:ILE:HD13	4:F:240:LEU:HB2	1.99	0.45
1:A:189:LEU:HD13	1:A:224:LEU:HD23	1.98	0.45
3:B:17:ASN:OD1	3:B:17:ASN:N	2.50	0.45
4:H:249:MET:HE2	4:H:249:MET:HB3	1.81	0.45
4:E:98:LYS:NZ	4:E:208:ASN:OD1	2.47	0.44
2:J:272:ARG:O	2:J:277:ARG:NH1	2.51	0.44
6:T:9:DA:H1'	6:T:10:DA:C8	2.51	0.44
4:F:121:LYS:HA	4:F:121:LYS:HD2	1.79	0.44
3:B:93:THR:HG23	3:B:96:TYR:HD2	1.81	0.44
2:J:97:MET:HE2	2:J:97:MET:HB2	1.74	0.44
2:J:173:VAL:HG13	2:J:177:LEU:HD23	1.99	0.44
2:J:87:LYS:NZ	2:J:200:SER:O	2.44	0.44
3:B:38:LYS:HE3	3:B:62:SER:HB2	1.99	0.44
1:A:5:TYR:HB3	1:A:109:LEU:HD11	1.99	0.44
2:J:30:SER:O	2:J:30:SER:OG	2.34	0.44
2:J:229:ASP:OD1	2:J:230:LYS:N	2.46	0.44
2:J:318:LYS:HE2	2:J:318:LYS:HB2	1.71	0.44
4:D:34:GLN:NE2	4:D:46:PRO:O	2.35	0.44
4:D:157:ILE:HG22	4:D:216:LYS:HB3	1.99	0.44
4:I:271:TYR:HB3	4:I:273:PHE:CE2	2.53	0.44
4:I:35:THR:OG1	4:I:36:ASP:N	2.51	0.44
4:I:196:ILE:HD13	4:I:207:LEU:HD23	1.98	0.44
4:I:259:ARG:NH2	4:I:278:GLU:OE2	2.51	0.44
4:D:199:THR:HG21	4:D:207:LEU:HB2	1.99	0.44
3:B:4:GLN:HB2	3:B:68:LEU:HD13	2.00	0.43
4:G:249:MET:HE3	4:G:249:MET:HB2	1.93	0.43
1:A:214:ASP:HB3	2:J:159:HIS:CD2	2.53	0.43
4:I:61:GLU:OE1	4:H:15:ARG:NH2	2.40	0.43
4:E:121:LYS:HA	4:E:121:LYS:HD2	1.80	0.43
3:B:94:ASP:OD1	3:B:94:ASP:N	2.50	0.43
3:B:127:LEU:O	3:B:130:HIS:ND1	2.44	0.43
5:C:51:G:O2'	5:C:55:A:N6	2.51	0.43
1:A:7:LEU:HD13	1:A:109:LEU:HD13	2.00	0.43
3:B:117:LYS:HA	3:B:117:LYS:HD2	1.72	0.43
4:D:99:VAL:HG11	4:D:196:ILE:HD11	2.00	0.43
2:J:151:PHE:HD2	5:C:1:U:H2'	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:E:279:ASN:ND2	4:F:59:GLY:O	2.52	0.43
1:A:207:GLY:HA2	4:I:17:GLU:HG3	2.00	0.43
3:B:153:GLN:OE1	3:B:154:SER:N	2.40	0.43
4:E:63:PRO:HA	4:E:64:PRO:HD3	1.92	0.43
4:H:166:ASP:OD1	4:H:166:ASP:N	2.52	0.43
6:T:8:DC:C4	6:T:9:DA:C6	3.06	0.43
6:T:17:DT:C2	4:D:328:MET:HE1	2.54	0.43
4:E:69:LYS:HG3	4:E:70:GLU:HG3	2.01	0.43
4:H:14:SER:OG	4:H:15:ARG:N	2.52	0.43
1:A:133:ARG:NH1	5:C:5:G:OP1	2.52	0.42
2:J:239:ILE:HD12	2:J:321:ILE:HD12	2.00	0.42
5:C:51:G:H22	5:C:55:A:P	2.42	0.42
3:B:20:ILE:HA	3:B:24:ILE:HD12	2.02	0.42
4:I:262:ASP:HB3	4:I:275:ILE:HG13	2.02	0.42
2:J:234:LEU:HD23	2:J:234:LEU:HA	1.92	0.42
2:J:326:GLU:O	2:J:330:ARG:HG2	2.19	0.42
4:F:236:LYS:HE2	4:F:236:LYS:HB2	1.84	0.42
4:G:35:THR:OG1	4:G:36:ASP:N	2.53	0.42
4:E:132:LEU:HG	4:E:136:LYS:HD2	2.01	0.42
4:F:331:LYS:HD2	4:F:331:LYS:HA	1.81	0.42
3:B:43:PHE:HE1	3:B:58:ILE:HD13	1.84	0.42
4:I:304:MET:HE3	4:I:311:LEU:HD12	2.02	0.42
6:T:48:DT:H2"	6:T:49:A:C8	2.55	0.42
4:D:35:THR:OG1	4:D:36:ASP:N	2.53	0.42
4:F:332:LYS:O	4:F:332:LYS:NZ	2.34	0.42
4:H:331:LYS:HA	4:H:331:LYS:HD3	1.83	0.42
2:J:188:ARG:HB2	2:J:208:TRP:CD2	2.55	0.42
3:B:3:SER:HA	3:B:60:VAL:O	2.18	0.42
4:I:284:ARG:HE	4:I:284:ARG:HB3	1.63	0.42
6:T:26:C:H4'	4:G:73:ASP:HB3	2.01	0.42
4:E:71:ILE:HG23	4:E:72:MET:SD	2.60	0.42
4:F:16:PRO:HD3	4:F:114:TYR:CD1	2.55	0.42
4:G:111:ASP:OD1	4:G:112:TYR:N	2.53	0.42
2:J:124:LYS:HD2	2:J:124:LYS:HA	1.76	0.42
4:G:39:LYS:HB2	4:G:39:LYS:HE3	1.87	0.42
4:H:103:LYS:NZ	4:H:123:ASP:OD1	2.43	0.42
1:A:22:ILE:HD12	1:A:23:ILE:HG12	2.02	0.41
2:J:114:LEU:O	2:J:118:MET:HG2	2.19	0.41
2:J:259:ILE:HD13	2:J:259:ILE:HA	1.89	0.41
3:B:2:PHE:HB3	3:B:90:ILE:HG22	2.01	0.41
3:B:145:LYS:HA	3:B:145:LYS:HD3	1.79	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:205:GLU:HG2	4:H:206:TYR:HD1	1.84	0.41
4:H:265:TYR:OH	4:H:273:PHE:O	2.29	0.41
7:N:12:DT:H2'	7:N:13:DT:H71	2.01	0.41
2:J:188:ARG:HB2	2:J:208:TRP:CG	2.55	0.41
3:B:153:GLN:CD	3:B:154:SER:H	2.27	0.41
6:T:10:DA:H3'	6:T:11:DG:C8	2.55	0.41
2:J:22:LYS:HD3	2:J:22:LYS:HA	1.77	0.41
6:T:32:DT:H1'	4:H:74:PRO:HG2	2.02	0.41
4:G:29:GLU:OE1	4:G:251:SER:N	2.54	0.41
4:G:125:TYR:OH	4:G:317:HIS:ND1	2.50	0.41
1:A:18:ALA:HB2	1:A:102:ILE:HB	2.02	0.41
5:C:36:A:H2'	5:C:37:A:C8	2.55	0.41
6:T:33:A:OP1	4:H:58:LYS:NZ	2.50	0.41
4:D:148:LEU:O	4:D:151:ASN:ND2	2.49	0.41
4:G:116:THR:O	4:G:119:TYR:HB2	2.21	0.41
4:F:250:HIS:HB3	4:F:252:GLN:OE1	2.21	0.41
1:A:153:ASP:OD1	1:A:153:ASP:N	2.54	0.41
2:J:316:SER:O	2:J:320:MET:HE1	2.20	0.41
4:E:14:SER:OG	4:E:15:ARG:N	2.54	0.41
4:E:143:ALA:O	4:E:174:SER:OG	2.31	0.41
4:F:166:ASP:OD1	4:F:166:ASP:N	2.51	0.41
4:F:252:GLN:NE2	4:G:54:VAL:HG12	2.36	0.41
4:G:89:ASP:OD1	4:G:89:ASP:N	2.54	0.41
1:A:43:LEU:HD23	1:A:43:LEU:HA	1.83	0.41
2:J:18:GLN:O	2:J:22:LYS:HG2	2.21	0.41
2:J:237:ILE:HG23	2:J:242:LEU:HD21	2.02	0.41
2:J:268:GLU:O	2:J:270:ASN:ND2	2.54	0.41
3:B:115:MET:HE2	3:B:115:MET:HB3	1.99	0.41
4:I:25:LEU:HD12	4:I:25:LEU:HA	1.96	0.41
5:C:35:U:O2'	4:D:327:GLY:O	2.39	0.41
4:D:300:LEU:HD23	4:D:315:ASP:HB3	2.03	0.41
4:F:271:TYR:HB3	4:F:273:PHE:CE2	2.56	0.41
4:G:106:MET:HE2	4:G:106:MET:HB3	1.86	0.41
4:G:278:GLU:OE1	4:G:281:GLY:N	2.54	0.41
4:H:275:ILE:HD13	4:H:290:PHE:HD2	1.86	0.41
1:A:247:GLU:H	1:A:247:GLU:HG2	1.66	0.41
2:J:10:ILE:HD11	2:J:23:TYR:HD1	1.85	0.40
2:J:133:LEU:HD23	2:J:133:LEU:HA	1.86	0.40
3:B:16:PRO:HA	3:B:19:ILE:HD12	2.03	0.40
4:G:121:LYS:HA	4:G:121:LYS:HD2	1.90	0.40
1:A:24:VAL:HG23	2:J:211:VAL:HG11	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:138:MET:SD	2:J:138:MET:N	2.94	0.40
3:B:123:LEU:HD13	5:C:54:U:N3	2.36	0.40
6:T:8:DC:H6	6:T:8:DC:H2'	1.72	0.40
4:G:313:ILE:HD12	4:G:313:ILE:HA	1.94	0.40
1:A:73:GLN:H	1:A:73:GLN:CD	2.28	0.40
2:J:177:LEU:HD21	2:J:212:PHE:CZ	2.57	0.40
2:J:320:MET:SD	2:J:320:MET:N	2.94	0.40
4:I:143:ALA:O	4:I:174:SER:OG	2.30	0.40
4:D:137:ARG:NH1	4:D:261:ILE:O	2.48	0.40
1:A:125:ILE:O	1:A:129:MET:HB2	2.22	0.40
1:A:221:ILE:HG22	2:J:68:ASP:HB3	2.02	0.40
2:J:97:MET:O	2:J:98:SER:C	2.65	0.40
5:C:17:A:O2'	4:G:21:PHE:O	2.35	0.40
2:J:303:VAL:HG13	2:J:320:MET:HG3	2.02	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	253/255 (99%)	236 (93%)	17 (7%)	0	100	100
2	J	338/344 (98%)	316 (94%)	22 (6%)	0	100	100
3	B	173/181 (96%)	153 (88%)	20 (12%)	0	100	100
4	D	291/335 (87%)	279 (96%)	12 (4%)	0	100	100
4	E	320/335 (96%)	305 (95%)	15 (5%)	0	100	100
4	F	321/335 (96%)	304 (95%)	17 (5%)	0	100	100
4	G	321/335 (96%)	308 (96%)	13 (4%)	0	100	100
4	H	321/335 (96%)	311 (97%)	10 (3%)	0	100	100
4	I	322/335 (96%)	314 (98%)	8 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	2660/2790 (95%)	2526 (95%)	134 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	A	219/219 (100%)	219 (100%)	0	100	100
2	J	299/304 (98%)	294 (98%)	5 (2%)	53	69
3	B	156/160 (98%)	155 (99%)	1 (1%)	78	80
4	D	265/299 (89%)	265 (100%)	0	100	100
4	E	289/299 (97%)	289 (100%)	0	100	100
4	F	290/299 (97%)	290 (100%)	0	100	100
4	G	290/299 (97%)	289 (100%)	1 (0%)	86	84
4	H	290/299 (97%)	288 (99%)	2 (1%)	76	78
4	I	291/299 (97%)	290 (100%)	1 (0%)	86	84
All	All	2389/2477 (96%)	2379 (100%)	10 (0%)	81	82

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	J	96	GLU
2	J	97	MET
2	J	98	SER
2	J	123	ASP
2	J	187	LEU
3	B	7	ILE
4	I	284	ARG
4	G	109	CYS
4	H	229	MET
4	H	279	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	ASN
1	A	237	ASN
2	J	4	ASN
2	J	12	GLN
2	J	238	ASN
2	J	311	GLN
4	D	144	ASN
4	D	194	GLN
4	D	250	HIS
4	E	41	HIS
4	G	77	GLN
4	H	186	ASN
4	H	279	ASN
4	H	308	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
5	C	59/60 (98%)	24 (40%)	0
6	T	0/48	-	-
7	N	0/11	-	-
All	All	59/119 (49%)	24 (40%)	0

All (24) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
5	C	3	U
5	C	9	G
5	C	15	U
5	C	16	C
5	C	19	U
5	C	20	U
5	C	21	A
5	C	22	A
5	C	25	A
5	C	26	G
5	C	27	G
5	C	32	U
5	C	33	G
5	C	34	U

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Mol	Chain	Res	Type
5	C	35	U
5	C	39	A
5	C	40	A
5	C	42	U
5	C	43	G
5	C	44	U
5	C	48	G
5	C	49	C
5	C	51	G
5	C	53	A

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](#)

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-39200. 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.