



Full wwPDB EM Validation Report ⓘ

Nov 21, 2022 – 01:27 AM EST

PDB ID : 7RB2
EMDB ID : EMD-24392
Title : Cryo-EM structure of SARS-CoV-2 NSP15 NendoU in BIS-Tris pH 6.0
Authors : Godoy, A.S.; Song, Y.; Nakamura, A.M.; Noske, G.D.; Gawriljuk, V.O.; Fernandes, R.S.; Oliva, G.
Deposited on : 2021-07-05
Resolution : 3.27 Å(reported)
Based on initial model : 7KEG

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.dev43
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : **FAILED**
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.3

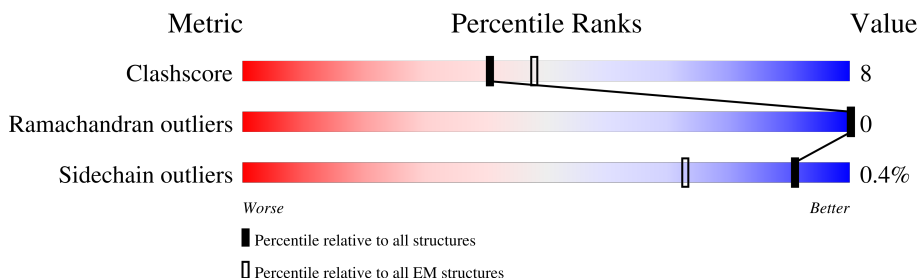
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.27 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	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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	349	80% 19% ..
1	B	349	80% 19% .
1	C	349	82% 17% ..
1	D	349	82% 17% .
1	E	349	79% 20% .
1	F	349	74% 25% ..

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 16386 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 Uridylate-specific endoribonuclease.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	347	Total	C	N	O	S	1	0
			2737	1760	444	522	11		
1	B	347	Total	C	N	O	S	1	0
			2737	1760	444	522	11		
1	C	347	Total	C	N	O	S	1	0
			2721	1750	442	518	11		
1	D	347	Total	C	N	O	S	1	0
			2725	1752	442	520	11		
1	E	347	Total	C	N	O	S	1	0
			2733	1757	443	522	11		
1	F	347	Total	C	N	O	S	1	0
			2733	1758	444	520	11		

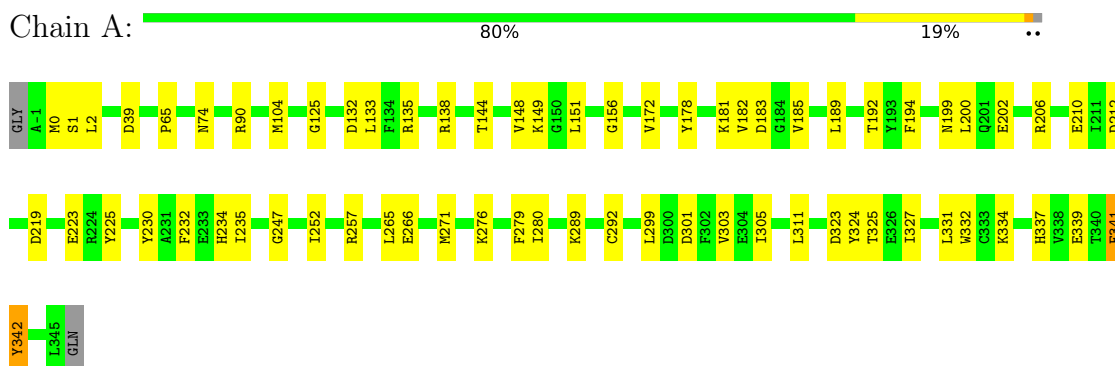
There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP P0DTD1
A	-1	ALA	-	expression tag	UNP P0DTD1
A	0	MET	-	expression tag	UNP P0DTD1
B	-2	GLY	-	expression tag	UNP P0DTD1
B	-1	ALA	-	expression tag	UNP P0DTD1
B	0	MET	-	expression tag	UNP P0DTD1
C	-2	GLY	-	expression tag	UNP P0DTD1
C	-1	ALA	-	expression tag	UNP P0DTD1
C	0	MET	-	expression tag	UNP P0DTD1
D	-2	GLY	-	expression tag	UNP P0DTD1
D	-1	ALA	-	expression tag	UNP P0DTD1
D	0	MET	-	expression tag	UNP P0DTD1
E	-2	GLY	-	expression tag	UNP P0DTD1
E	-1	ALA	-	expression tag	UNP P0DTD1
E	0	MET	-	expression tag	UNP P0DTD1
F	-2	GLY	-	expression tag	UNP P0DTD1
F	-1	ALA	-	expression tag	UNP P0DTD1
F	0	MET	-	expression tag	UNP P0DTD1

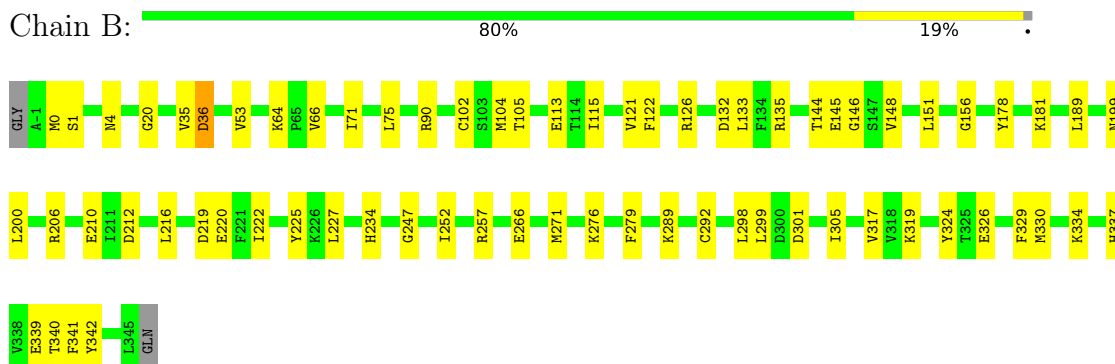
3 Residue-property plots

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.

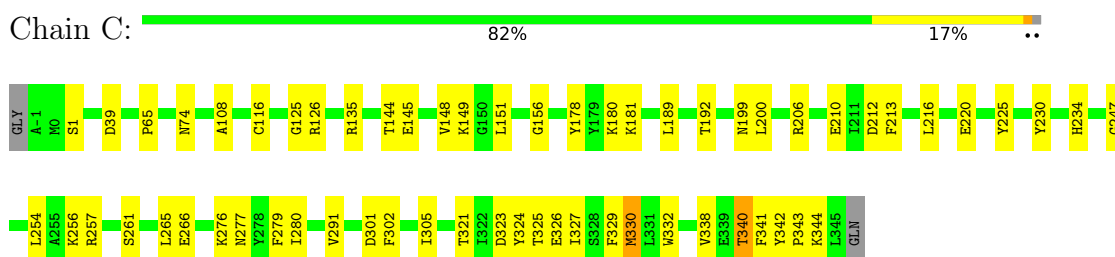
- Molecule 1: Uridylate-specific endoribonuclease




- Molecule 1: Uridylate-specific endoribonuclease

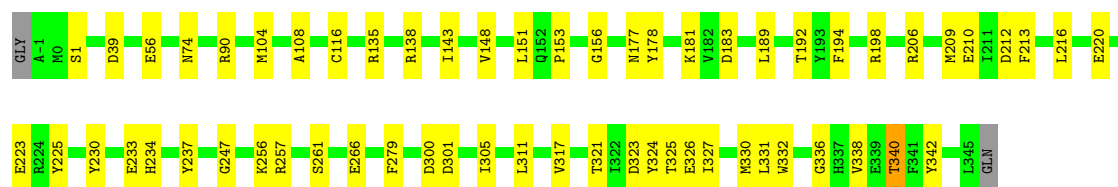


- Molecule 1: Uridylate-specific endoribonuclease




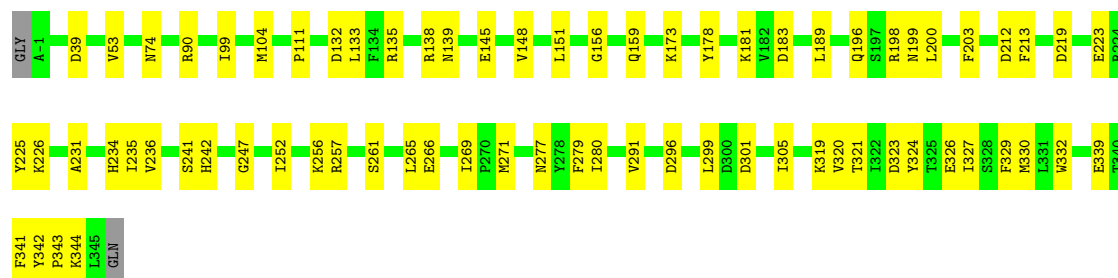
- Molecule 1: Uridylate-specific endoribonuclease

Chain D:  82% 17%



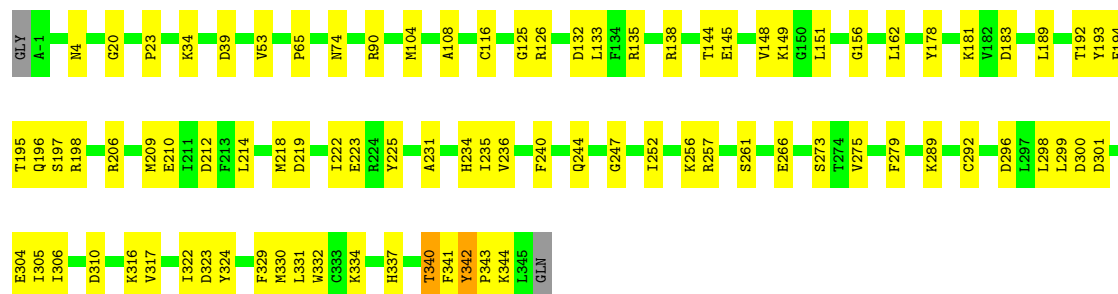
• Molecule 1: Uridylate-specific endoribonuclease

Chain E:  79% 20%



• Molecule 1: Uridylate-specific endoribonuclease

Chain F:  74% 25%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	263322	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 ($e^-/\text{\AA}^2$)	43.254543	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor

5 Model quality

5.1 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 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	0.41	0/2797	0.50	0/3796
1	B	0.41	0/2797	0.49	0/3796
1	C	0.41	0/2781	0.49	0/3778
1	D	0.41	0/2785	0.49	0/3783
1	E	0.42	0/2793	0.50	0/3792
1	F	0.42	0/2793	0.50	0/3791
All	All	0.41	0/16746	0.50	0/22736

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

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	2737	0	2726	42	0
1	B	2737	0	2726	38	0
1	C	2721	0	2696	36	0
1	D	2725	0	2700	38	0
1	E	2733	0	2715	42	0
1	F	2733	0	2722	56	0
All	All	16386	0	16285	245	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 (245) 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:F:65:PRO:HG3	1:F:162:LEU:HD12	1.56	0.87
1:E:266:GLU:HB2	1:E:279:PHE:HB3	1.66	0.78
1:D:266:GLU:HB2	1:D:279:PHE:HB3	1.67	0.76
1:B:266:GLU:HB2	1:B:279:PHE:HB3	1.68	0.76
1:E:257:ARG:HH12	1:E:261:SER:HB3	1.51	0.75
1:F:266:GLU:HB3	1:F:279:PHE:HB3	1.68	0.74
1:E:235:ILE:HG13	1:E:236:VAL:HG23	1.69	0.73
1:F:235:ILE:HG13	1:F:236:VAL:HG23	1.70	0.73
1:C:266:GLU:HB2	1:C:279:PHE:HB3	1.69	0.72
1:F:212:ASP:OD1	1:F:225:TYR:OH	2.08	0.71
1:E:135:ARG:O	1:E:181:LYS:NZ	2.24	0.71
1:F:193:TYR:OH	1:F:304:GLU:OE1	2.08	0.71
1:D:257:ARG:HH12	1:D:261:SER:HB3	1.55	0.70
1:D:301:ASP:O	1:D:305:ILE:HG13	1.92	0.69
1:A:342:TYR:HD2	1:A:342:TYR:C	1.96	0.68
1:E:213:PHE:O	1:E:256:LYS:NZ	2.26	0.68
1:E:132:ASP:OD2	1:E:133:LEU:N	2.26	0.68
1:B:102:CYS:SG	1:B:105:THR:OG1	2.49	0.68
1:F:332:TRP:HB2	1:F:340:THR:HG22	1.75	0.68
1:A:266:GLU:HB2	1:A:279:PHE:HB3	1.75	0.68
1:F:132:ASP:OD1	1:F:133:LEU:N	2.27	0.68
1:B:132:ASP:OD2	1:B:133:LEU:N	2.27	0.67
1:A:132:ASP:OD2	1:A:133:LEU:N	2.27	0.67
1:A:135:ARG:O	1:A:181:LYS:NZ	2.28	0.67
1:F:135:ARG:O	1:F:181:LYS:NZ	2.27	0.67
1:F:214:LEU:O	1:F:256:LYS:NZ	2.28	0.66
1:E:212:ASP:OD2	1:E:225:TYR:OH	2.07	0.66
1:A:138:ARG:HE	1:A:183:ASP:HA	1.61	0.66
1:B:135:ARG:O	1:B:181:LYS:NZ	2.30	0.65
1:D:135:ARG:O	1:D:181:LYS:NZ	2.29	0.65
1:E:265:LEU:HD22	1:E:280:ILE:HG12	1.79	0.64
1:C:257:ARG:HH12	1:C:261:SER:HB3	1.63	0.64
1:F:257:ARG:HH12	1:F:261:SER:HB3	1.64	0.63
1:D:213:PHE:O	1:D:256:LYS:NZ	2.32	0.63
1:A:342:TYR:C	1:A:342:TYR:CD2	2.71	0.62
1:B:301:ASP:O	1:B:305:ILE:HG13	2.00	0.61
1:D:330:MET:HG2	1:D:342:TYR:CZ	2.36	0.61
1:D:332:TRP:HB2	1:D:340:THR:HG22	1.81	0.61
1:E:301:ASP:O	1:E:305:ILE:HG13	2.00	0.61
1:D:212:ASP:OD2	1:D:225:TYR:OH	2.11	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:301:ASP:O	1:A:305:ILE:HG13	2.02	0.60
1:B:334:LYS:O	1:B:337:HIS:ND1	2.36	0.59
1:C:213:PHE:O	1:C:256:LYS:NZ	2.32	0.58
1:A:212:ASP:OD2	1:A:225:TYR:OH	2.11	0.58
1:F:334:LYS:O	1:F:337:HIS:ND1	2.34	0.58
1:A:342:TYR:CD2	1:A:342:TYR:O	2.57	0.57
1:B:126:ARG:HG2	1:B:145:GLU:HG2	1.85	0.57
1:C:265:LEU:HD22	1:C:280:ILE:HG12	1.86	0.57
1:C:39:ASP:O	1:F:90:ARG:NH1	2.38	0.56
1:F:244:GLN:NE2	1:F:289:LYS:HG3	2.20	0.56
1:D:330:MET:HG2	1:D:342:TYR:OH	2.06	0.56
1:B:257:ARG:HD2	1:B:257:ARG:O	2.06	0.55
1:E:198:ARG:HH21	1:E:203:PHE:HZ	1.54	0.55
1:F:301:ASP:O	1:F:305:ILE:HG12	2.07	0.55
1:F:317:VAL:HG22	1:F:330:MET:HG3	1.88	0.55
1:B:212:ASP:OD2	1:B:225:TYR:OH	2.12	0.54
1:F:138:ARG:NH1	1:F:183:ASP:HA	2.22	0.54
1:D:148:VAL:HG13	1:D:151:LEU:HB2	1.89	0.54
1:C:332:TRP:HB2	1:C:340:THR:HG22	1.89	0.54
1:A:271:MET:O	1:A:276:LYS:NZ	2.38	0.53
1:C:74:ASN:ND2	1:C:323:ASP:O	2.40	0.53
1:C:330:MET:SD	1:C:344:LYS:NZ	2.68	0.53
1:D:198:ARG:HD3	1:D:206:ARG:HH11	1.73	0.53
1:A:342:TYR:HD2	1:A:342:TYR:O	1.91	0.53
1:A:90:ARG:NH1	1:E:39:ASP:O	2.41	0.53
1:B:317:VAL:HG22	1:B:330:MET:SD	2.50	0.53
1:B:271:MET:O	1:B:276:LYS:NZ	2.40	0.52
1:A:148:VAL:HG13	1:A:151:LEU:HB2	1.91	0.52
1:D:209:MET:HE3	1:D:300:ASP:OD1	2.10	0.52
1:A:332:TRP:O	1:A:339:GLU:N	2.39	0.52
1:B:189:LEU:HD12	1:B:324:TYR:HD1	1.75	0.52
1:C:135:ARG:O	1:C:181:LYS:NZ	2.42	0.51
1:F:273:SER:OG	1:F:275:VAL:O	2.25	0.51
1:F:195:THR:HG22	1:F:197:SER:H	1.76	0.51
1:A:265:LEU:HD22	1:A:280:ILE:HG12	1.93	0.51
1:A:199:ASN:OD1	1:A:200:LEU:N	2.44	0.51
1:A:234:HIS:O	1:A:247:GLY:HA3	2.10	0.51
1:E:330:MET:SD	1:E:344:LYS:HD3	2.51	0.51
1:C:330:MET:SD	1:C:342:TYR:HE2	2.34	0.50
1:D:230:TYR:HE1	1:D:311:LEU:HD11	1.74	0.50
1:F:193:TYR:C	1:F:322:ILE:HD12	2.32	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:125:GLY:HA3	1:C:144:THR:HG22	1.93	0.50
1:A:334:LYS:O	1:A:337:HIS:ND1	2.46	0.49
1:D:230:TYR:HA	1:D:338:VAL:HB	1.93	0.49
1:F:330:MET:SD	1:F:344:LYS:HD3	2.53	0.49
1:D:56:GLU:OE1	1:D:104:MET:HE1	2.13	0.49
1:D:189:LEU:HD12	1:D:324:TYR:HD2	1.77	0.49
1:F:289:LYS:HB3	1:F:292:CYS:SG	2.53	0.49
1:F:209:MET:HE3	1:F:300:ASP:OD1	2.13	0.49
1:F:231:ALA:O	1:F:235:ILE:HG12	2.13	0.49
1:B:252:ILE:HD13	1:B:299:LEU:HD22	1.95	0.49
1:B:199:ASN:OD1	1:B:200:LEU:N	2.46	0.48
1:E:138:ARG:NH1	1:E:183:ASP:HA	2.28	0.48
1:A:39:ASP:O	1:D:90:ARG:NH1	2.47	0.48
1:D:138:ARG:NH1	1:D:183:ASP:HA	2.29	0.48
1:A:202:GLU:OE1	1:A:202:GLU:N	2.46	0.48
1:A:252:ILE:HD13	1:A:299:LEU:HD22	1.95	0.48
1:D:206:ARG:N	1:D:210:GLU:OE2	2.26	0.48
1:A:289:LYS:HB3	1:A:292:CYS:SG	2.53	0.48
1:F:198:ARG:HD3	1:F:206:ARG:HH11	1.78	0.48
1:F:330:MET:HB3	1:F:342:TYR:CE2	2.49	0.48
1:D:234:HIS:O	1:D:247:GLY:HA3	2.14	0.47
1:E:148:VAL:HG13	1:E:151:LEU:HB2	1.96	0.47
1:E:321:THR:OG1	1:E:326:GLU:OE1	2.26	0.47
1:C:199:ASN:OD1	1:C:200:LEU:N	2.47	0.47
1:E:219:ASP:O	1:E:223:GLU:OE1	2.31	0.47
1:E:223:GLU:O	1:E:226:LYS:HD2	2.14	0.47
1:A:232:PHE:HZ	1:A:303:VAL:HG12	1.79	0.47
1:F:330:MET:SD	1:F:342:TYR:HE2	2.36	0.47
1:B:90:ARG:NH1	1:F:39:ASP:O	2.46	0.47
1:D:74:ASN:ND2	1:D:323:ASP:O	2.43	0.47
1:D:39:ASP:O	1:E:90:ARG:NH1	2.47	0.47
1:E:269:ILE:HG22	1:E:271:MET:HG2	1.96	0.47
1:F:74:ASN:ND2	1:F:323:ASP:O	2.40	0.47
1:F:219:ASP:O	1:F:223:GLU:OE1	2.31	0.47
1:F:196:GLN:HB2	1:F:296:ASP:OD2	2.15	0.47
1:B:71:ILE:O	1:B:75:LEU:HD23	2.15	0.47
1:C:216:LEU:HB3	1:C:220:GLU:HB2	1.97	0.47
1:E:199:ASN:OD1	1:E:200:LEU:N	2.48	0.47
1:B:330:MET:HG2	1:B:342:TYR:OH	2.15	0.46
1:E:252:ILE:HD13	1:E:299:LEU:HD22	1.96	0.46
1:A:230:TYR:HE1	1:A:311:LEU:HD21	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:53:VAL:HG22	1:F:104:MET:SD	2.55	0.46
1:A:189:LEU:HD12	1:A:324:TYR:HD1	1.80	0.46
1:C:180:LYS:HB3	1:C:189:LEU:HD21	1.97	0.46
1:E:189:LEU:HD12	1:E:324:TYR:HD2	1.80	0.46
1:F:218:MET:O	1:F:222:ILE:HG13	2.15	0.46
1:F:189:LEU:HD12	1:F:324:TYR:HD1	1.81	0.46
1:A:156:GLY:HA3	1:A:178:TYR:CE1	2.51	0.46
1:C:148:VAL:HG13	1:C:151:LEU:HB2	1.98	0.46
1:A:74:ASN:ND2	1:A:323:ASP:O	2.41	0.46
1:C:234:HIS:O	1:C:247:GLY:HA3	2.15	0.46
1:B:289:LYS:HB3	1:B:292:CYS:SG	2.56	0.46
1:E:231:ALA:O	1:E:235:ILE:HG12	2.15	0.46
1:B:234:HIS:O	1:B:247:GLY:HA3	2.15	0.45
1:B:113:GLU:HG2	1:B:115:ILE:HG22	1.97	0.45
1:E:139:ASN:HA	1:E:181:LYS:O	2.16	0.45
1:B:148:VAL:HG13	1:B:151:LEU:HB2	1.97	0.45
1:B:206:ARG:N	1:B:210:GLU:OE1	2.26	0.45
1:D:311:LEU:HB3	1:D:336:GLY:O	2.16	0.45
1:F:310:ASP:HB3	1:F:316:LYS:NZ	2.32	0.45
1:C:329:PHE:HB3	1:C:341:PHE:CE1	2.51	0.45
1:D:108:ALA:HB3	1:D:116:CYS:SG	2.56	0.45
1:D:321:THR:OG1	1:D:326:GLU:OE2	2.22	0.45
1:C:189:LEU:HD12	1:C:324:TYR:HD1	1.81	0.45
1:D:325:THR:HG23	1:D:327:ILE:HD11	1.98	0.45
1:B:222:ILE:HG23	1:B:227:LEU:HB2	1.97	0.45
1:A:104:MET:HG2	1:F:23:PRO:HB3	1.98	0.45
1:F:206:ARG:N	1:F:210:GLU:OE1	2.25	0.45
1:A:135:ARG:HH22	1:A:149:LYS:HZ3	1.63	0.45
1:A:235:ILE:HD13	1:A:341:PHE:HB2	1.99	0.45
1:C:301:ASP:O	1:C:305:ILE:HG23	2.17	0.44
1:D:331:LEU:HD12	1:D:340:THR:O	2.17	0.44
1:F:210:GLU:OE2	1:F:298:LEU:HD12	2.17	0.44
1:B:64:LYS:O	1:B:66:VAL:HG13	2.17	0.44
1:E:332:TRP:O	1:E:339:GLU:N	2.31	0.44
1:C:206:ARG:N	1:C:210:GLU:OE2	2.25	0.44
1:E:53:VAL:HG22	1:E:104:MET:SD	2.58	0.44
1:F:240:PHE:CE1	1:F:257:ARG:HD3	2.52	0.44
1:D:143:ILE:O	1:D:143:ILE:HG13	2.18	0.44
1:F:148:VAL:HG13	1:F:151:LEU:HB2	2.00	0.44
1:D:317:VAL:HG22	1:D:330:MET:SD	2.57	0.44
1:C:325:THR:HG23	1:C:327:ILE:HD11	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:153:PRO:HB2	1:D:177:ASN:HB3	1.98	0.44
1:E:320:VAL:O	1:E:327:ILE:N	2.35	0.44
1:F:108:ALA:HB3	1:F:116:CYS:SG	2.58	0.44
1:F:192:THR:HG23	1:F:194:PHE:CE1	2.53	0.44
1:B:330:MET:HB3	1:B:342:TYR:CE1	2.52	0.43
1:E:138:ARG:HH12	1:E:183:ASP:HA	1.83	0.43
1:C:149:LYS:HD3	1:C:149:LYS:C	2.38	0.43
1:B:144:THR:HG22	1:B:146:GLY:H	1.83	0.43
1:F:316:LYS:HZ3	1:F:331:LEU:HD23	1.83	0.43
1:A:219:ASP:O	1:A:223:GLU:HG2	2.18	0.43
1:B:329:PHE:HB3	1:B:341:PHE:CE1	2.53	0.43
1:C:230:TYR:HA	1:C:338:VAL:HB	2.01	0.43
1:F:34:LYS:NZ	1:F:39:ASP:OD2	2.49	0.43
1:A:206:ARG:N	1:A:210:GLU:OE1	2.27	0.43
1:F:234:HIS:O	1:F:247:GLY:HA3	2.18	0.43
1:F:252:ILE:HD13	1:F:299:LEU:HD22	2.00	0.43
1:A:125:GLY:HA3	1:A:144:THR:HG22	2.00	0.43
1:D:220:GLU:O	1:D:223:GLU:HG3	2.19	0.43
1:A:257:ARG:HD2	1:A:257:ARG:O	2.19	0.43
1:F:149:LYS:HD3	1:F:149:LYS:C	2.39	0.42
1:F:193:TYR:CE2	1:F:305:ILE:HD11	2.54	0.42
1:A:192:THR:HG23	1:A:194:PHE:CE1	2.54	0.42
1:C:321:THR:OG1	1:C:326:GLU:OE1	2.31	0.42
1:E:196:GLN:HB2	1:E:296:ASP:OD2	2.20	0.42
1:E:74:ASN:ND2	1:E:323:ASP:O	2.46	0.42
1:C:156:GLY:HA3	1:C:178:TYR:CE1	2.54	0.42
1:D:233:GLU:HA	1:D:237:TYR:HB2	2.01	0.42
1:F:126:ARG:HG2	1:F:145:GLU:HG2	2.01	0.42
1:B:4:ASN:HA	1:B:20:GLY:O	2.20	0.42
1:E:319:LYS:HE3	1:E:326:GLU:HB3	2.02	0.42
1:E:329:PHE:HB3	1:E:341:PHE:CE1	2.55	0.42
1:B:339:GLU:HG2	1:B:340:THR:HG23	2.02	0.42
1:E:99:ILE:HD13	1:E:111:PRO:HA	2.01	0.42
1:B:53:VAL:HG22	1:B:104:MET:SD	2.60	0.42
1:B:219:ASP:O	1:B:220:GLU:C	2.58	0.42
1:C:212:ASP:OD2	1:C:225:TYR:OH	2.15	0.42
1:F:125:GLY:HA3	1:F:144:THR:HG22	2.01	0.42
1:E:145:GLU:OE2	1:E:173:LYS:HE3	2.20	0.42
1:E:156:GLY:HA3	1:E:178:TYR:CE1	2.55	0.42
1:F:193:TYR:O	1:F:194:PHE:HD1	2.03	0.42
1:B:210:GLU:OE2	1:B:298:LEU:HD12	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:126:ARG:HG2	1:C:145:GLU:HG2	2.02	0.41
1:F:4:ASN:HA	1:F:20:GLY:O	2.20	0.41
1:C:192:THR:HB	1:C:321:THR:HG22	2.02	0.41
1:C:276:LYS:HE3	1:C:276:LYS:HB3	1.90	0.41
1:F:289:LYS:HD3	1:F:292:CYS:SG	2.60	0.41
1:A:0:MET:SD	1:A:1:SER:N	2.93	0.41
1:D:330:MET:HB3	1:D:342:TYR:CE1	2.55	0.41
1:E:234:HIS:O	1:E:247:GLY:HA3	2.20	0.41
1:E:319:LYS:HE3	1:E:326:GLU:CB	2.50	0.41
1:A:325:THR:HG23	1:A:327:ILE:HD11	2.02	0.41
1:B:121:VAL:HG12	1:B:122:PHE:N	2.36	0.41
1:F:329:PHE:CE2	1:F:343:PRO:HG3	2.56	0.41
1:A:65:PRO:HG2	1:A:172:VAL:HG21	2.03	0.41
1:B:0:MET:SD	1:B:1:SER:N	2.94	0.41
1:C:277:ASN:HB3	1:C:291:VAL:HG23	2.03	0.41
1:E:329:PHE:CE1	1:E:343:PRO:HG3	2.56	0.41
1:F:329:PHE:HB3	1:F:341:PHE:CE1	2.56	0.41
1:B:35:VAL:HG12	1:B:36[B]:ASP:OD1	2.21	0.41
1:A:331:LEU:HB2	1:A:341:PHE:CE1	2.55	0.41
1:A:1:SER:OG	1:A:2:LEU:N	2.54	0.41
1:A:182:VAL:O	1:A:185:VAL:HG12	2.20	0.41
1:C:108:ALA:HB3	1:C:116:CYS:SG	2.60	0.41
1:C:200:LEU:HD21	1:C:254:LEU:HB3	2.03	0.41
1:B:319:LYS:HE3	1:B:326:GLU:HB3	2.03	0.41
1:D:192:THR:HG23	1:D:194:PHE:CE2	2.55	0.40
1:E:241:SER:OG	1:E:242:HIS:ND1	2.54	0.40
1:D:189:LEU:HD12	1:D:324:TYR:CD2	2.56	0.40
1:E:277:ASN:HB3	1:E:291:VAL:HG23	2.02	0.40
1:E:342:TYR:HA	1:E:343:PRO:HD3	1.94	0.40
1:F:156:GLY:HA3	1:F:178:TYR:HE1	1.86	0.40
1:E:159:GLN:OE1	1:E:159:GLN:N	2.54	0.40
1:F:306:ILE:HD12	1:F:306:ILE:HA	1.94	0.40
1:B:216:LEU:HB3	1:B:220:GLU:HB2	2.03	0.40
1:C:65:PRO:O	1:C:126:ARG:NH1	2.44	0.40
1:C:302:PHE:HA	1:C:305:ILE:HG12	2.04	0.40
1:C:342:TYR:HA	1:C:343:PRO:HD3	1.96	0.40
1:D:156:GLY:HA3	1:D:178:TYR:CE1	2.55	0.40
1:A:156:GLY:HA3	1:A:178:TYR:HE1	1.86	0.40
1:B:156:GLY:HA3	1:B:178:TYR:CE1	2.56	0.40
1:C:1:SER:HB2	1:D:1:SER:HB2	2.04	0.40
1:D:212:ASP:O	1:D:216:LEU:HB2	2.21	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	346/349 (99%)	327 (94%)	19 (6%)	0	100	100
1	B	346/349 (99%)	327 (94%)	19 (6%)	0	100	100
1	C	346/349 (99%)	330 (95%)	16 (5%)	0	100	100
1	D	346/349 (99%)	327 (94%)	19 (6%)	0	100	100
1	E	346/349 (99%)	331 (96%)	15 (4%)	0	100	100
1	F	346/349 (99%)	328 (95%)	18 (5%)	0	100	100
All	All	2076/2094 (99%)	1970 (95%)	106 (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	307/309 (99%)	305 (99%)	2 (1%)	84	90
1	B	307/309 (99%)	305 (99%)	2 (1%)	84	90
1	C	303/309 (98%)	301 (99%)	2 (1%)	84	90
1	D	304/309 (98%)	303 (100%)	1 (0%)	92	96
1	E	306/309 (99%)	306 (100%)	0	100	100
1	F	306/309 (99%)	304 (99%)	2 (1%)	84	90

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	1833/1854 (99%)	1824 (100%)	9 (0%)	91	93

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	341	PHE
1	A	342	TYR
1	B	36[A]	ASP
1	B	36[B]	ASP
1	C	330	MET
1	C	340	THR
1	D	340	THR
1	F	340	THR
1	F	342	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	337	HIS
1	B	249	HIS
1	E	249	HIS
1	F	244	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

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-24392. These allow visual inspection of the internal detail of the map and identification of artifacts.

No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

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 surface views

This section was not generated.

6.5 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.