



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 12:55 am BST

PDB ID : 6U8K
Title : Crystal structure of hepatitis C virus IRES junction IIIabc in complex with Fab HCV3
Authors : Koirala, D.; Lewicka, A.; Koldobskaya, Y.; Huang, H.; Piccirilli, J.A.
Deposited on : 2019-09-05
Resolution : 2.75 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

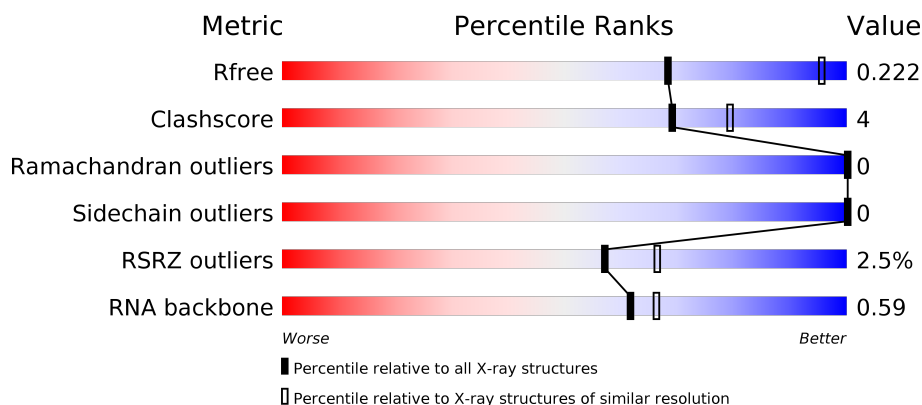
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION





The reported resolution of this entry is 2.75 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)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1235 (2.78-2.74)
Clashscore	141614	1277 (2.78-2.74)
Ramachandran outliers	138981	1257 (2.78-2.74)
Sidechain outliers	138945	1257 (2.78-2.74)
RSRZ outliers	127900	1207 (2.78-2.74)
RNA backbone	3102	1060 (3.02-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower 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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	68	 76% 22%
1	B	68	 79% 16%
1	C	68	 76% 21%
2	D	232	 3% 88% 7%

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Mol	Chain	Length	Quality of chain
2	F	232	<div><div></div><div>3%</div><div>88%</div><div>8%</div><div></div></div>
2	H	232	<div><div></div><div>3%</div><div>83%</div><div>13%</div><div></div></div>
3	E	215	<div><div></div><div>%</div><div>88%</div><div>12%</div><div></div></div>
3	G	215	<div><div></div><div>4%</div><div>91%</div><div>9%</div><div></div></div>
3	L	215	<div><div></div><div>2%</div><div>91%</div><div>9%</div><div></div></div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 14455 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, 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 RNA chain called JIIIabc RNA (68-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	68	Total	C	N	O	P	0	0	0
			1460	650	270	472	68			
1	B	68	Total	C	N	O	P	0	0	0
			1460	650	270	472	68			
1	C	68	Total	C	N	O	P	0	0	0
			1460	650	270	472	68			

- Molecule 2 is a protein called Heavy chain of Fab HCV3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	222	Total	C	N	O	S	0	0	0
			1661	1046	283	327	5			
2	F	222	Total	C	N	O	S	0	0	0
			1661	1046	283	327	5			
2	H	222	Total	C	N	O	S	0	0	0
			1661	1046	283	327	5			

- Molecule 3 is a protein called Light chain of Fab HCV3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	215	Total	C	N	O	S	0	0	0
			1644	1025	275	338	6			
3	G	215	Total	C	N	O	S	0	0	0
			1644	1025	275	338	6			
3	L	215	Total	C	N	O	S	0	0	0
			1644	1025	275	338	6			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	20	Total	O	0	0
			20	20		

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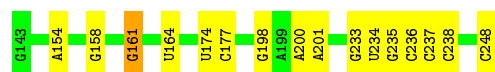
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	18	Total 18	O 18	0	0
4	C	5	Total 5	O 5	0	0
4	D	19	Total 19	O 19	0	0
4	E	16	Total 16	O 16	0	0
4	F	14	Total 14	O 14	0	0
4	G	12	Total 12	O 12	0	0
4	H	26	Total 26	O 26	0	0
4	L	30	Total 30	O 30	0	0

3 Residue-property plots [i](#)


These plots are drawn for all protein, RNA and DNA 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 and electron density. 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. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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: JIIIabc RNA (68-MER)

Chain A: 



- Molecule 1: JIIIabc RNA (68-MER)

Chain B: 




- Molecule 1: JIIIabc RNA (68-MER)

Chain C: 




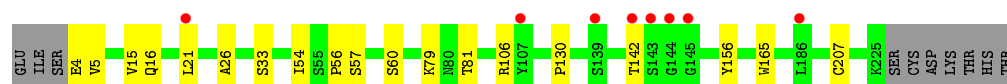
- Molecule 2: Heavy chain of Fab HCV3

Chain D: 

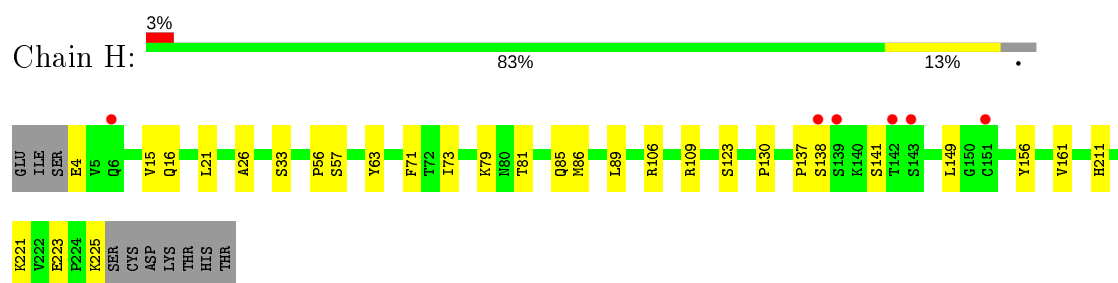


- Molecule 2: Heavy chain of Fab HCV3

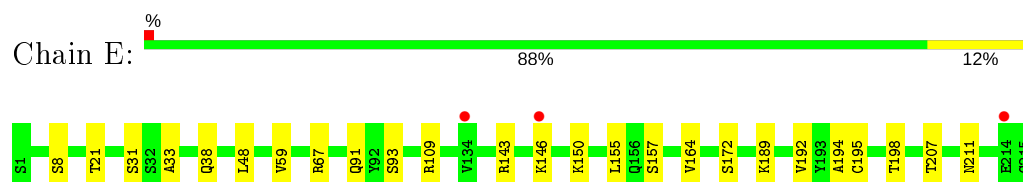
Chain F: 



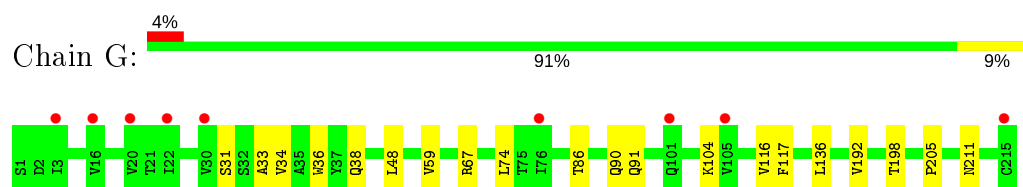
- Molecule 2: Heavy chain of Fab HCV3



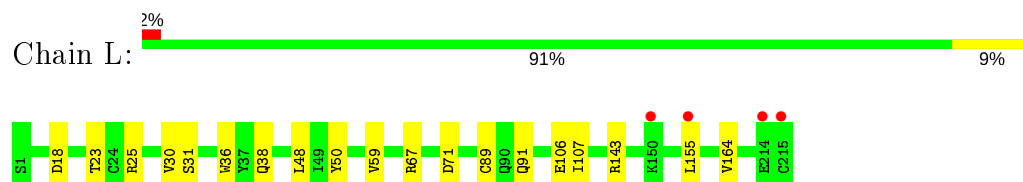
- Molecule 3: Light chain of Fab HCV3



- Molecule 3: Light chain of Fab HCV3



- Molecule 3: Light chain of Fab HCV3



4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	173.29 Å 173.29 Å 140.49 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	173.29 – 2.75 173.29 – 2.75	Depositor EDS
% Data completeness (in resolution range)	98.5 (173.29-2.75) 94.3 (173.29-2.75)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.95 (at 2.73 Å)	Xtriage
Refinement program	PHENIX (1.14_3260)	Depositor
R, R_{free}	0.186 , 0.222 0.187 , 0.222	Depositor DCC
R_{free} test set	2006 reflections (1.89%)	wwPDB-VP
Wilson B-factor (Å ²)	59.3	Xtriage
Anisotropy	0.163	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 64.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.177 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	14455	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.93% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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	0.54	0/1633	1.02	0/2546
1	B	0.55	0/1633	1.03	0/2546
1	C	0.45	0/1633	0.97	0/2546
2	D	0.45	0/1702	0.66	0/2317
2	F	0.43	0/1702	0.64	0/2317
2	H	0.46	0/1702	0.65	0/2317
3	E	0.48	0/1679	0.66	0/2279
3	G	0.44	0/1679	0.62	0/2279
3	L	0.48	0/1679	0.63	0/2279
All	All	0.48	0/15042	0.79	0/21426

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	1460	0	739	6	0
1	B	1460	0	739	5	0
1	C	1460	0	739	6	0
2	D	1661	0	1624	13	0
2	F	1661	0	1624	13	0
2	H	1661	0	1624	19	0
3	E	1644	0	1595	16	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	G	1644	0	1595	12	0
3	L	1644	0	1595	13	0
4	A	20	0	0	0	0
4	B	18	0	0	0	0
4	C	5	0	0	0	0
4	D	19	0	0	0	0
4	E	16	0	0	0	0
4	F	14	0	0	0	0
4	G	12	0	0	0	0
4	H	26	0	0	1	0
4	L	30	0	0	0	0
All	All	14455	0	11874	94	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 4.

All (94) 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)
3:G:198:THR:HG22	3:G:205:PRO:HB3	1.70	0.72
2:D:26:ALA:HA	2:D:81:THR:HG22	1.72	0.70
2:F:26:ALA:HA	2:F:81:THR:HG22	1.77	0.65
2:H:26:ALA:HA	2:H:81:THR:HG22	1.78	0.64
2:D:137:PRO:HG3	2:D:149:LEU:HB3	1.79	0.64
2:F:33:SER:HA	2:F:56:PRO:HB2	1.80	0.64
2:H:33:SER:HA	2:H:56:PRO:HB2	1.80	0.64
2:H:137:PRO:HG3	2:H:149:LEU:HB3	1.81	0.62
3:E:38:GLN:HB2	3:E:48:LEU:HD11	1.82	0.62
1:C:161:G:N7	2:F:106:ARG:NH1	2.48	0.61
2:F:33:SER:O	2:F:57:SER:HB3	2.01	0.61
2:D:33:SER:HA	2:D:56:PRO:HB2	1.82	0.61
3:E:21:THR:HG21	3:L:23:THR:HG21	1.85	0.59
3:G:31:SER:O	3:G:67:ARG:NH1	2.37	0.58
3:E:157:SER:HB2	3:L:155:LEU:HD21	1.85	0.58
2:D:4:GLU:HG2	2:D:5:VAL:H	1.69	0.57
1:B:161:G:N7	2:D:106:ARG:NH1	2.53	0.56
1:C:198:G:N2	1:C:201:A:OP2	2.40	0.55
2:F:130:PRO:HB3	2:F:156:TYR:HB3	1.90	0.54
2:H:33:SER:O	2:H:57:SER:HB3	2.07	0.54
2:D:4:GLU:HG2	2:D:5:VAL:N	2.24	0.53
2:H:4:GLU:OE1	2:H:4:GLU:N	2.42	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:150:LYS:HB2	3:E:194:ALA:HB3	1.91	0.52
3:G:38:GLN:HB2	3:G:48:LEU:HD11	1.92	0.52
3:L:30:VAL:HG11	3:L:91:GLN:HB2	1.90	0.52
3:L:38:GLN:HB2	3:L:48:LEU:HD11	1.91	0.52
2:H:15:VAL:HG11	2:H:89:LEU:HD13	1.91	0.51
1:C:148:U:H2'	1:C:149:C:C6	2.45	0.51
2:H:79:LYS:O	2:H:81:THR:HG23	2.11	0.51
1:C:201:A:C5	1:C:219:C:H1'	2.46	0.51
3:G:33:ALA:O	3:G:91:GLN:HA	2.11	0.50
2:H:161:VAL:HG12	2:H:211:HIS:CD2	2.46	0.50
3:L:48:LEU:HA	3:L:59:VAL:HG21	1.93	0.50
2:D:154:LYS:HA	2:D:188:SER:HB2	1.94	0.49
3:G:192:VAL:HG22	3:G:211:ASN:OD1	2.12	0.49
1:C:237:C:H2'	1:C:238:C:C6	2.48	0.49
3:E:146:LYS:HB3	3:E:198:THR:HG23	1.95	0.49
2:H:63:TYR:CZ	2:H:73:ILE:HG22	2.47	0.49
2:D:15:VAL:CG2	2:D:21:LEU:HD22	2.43	0.48
2:D:4:GLU:N	2:D:4:GLU:OE1	2.46	0.48
1:B:177:C:H2'	1:B:178:C:C6	2.47	0.48
1:C:237:C:H2'	1:C:238:C:H6	1.78	0.48
1:B:235:G:H1'	1:B:236:C:C6	2.49	0.47
3:E:189:LYS:HB3	3:E:189:LYS:HE2	1.67	0.47
2:H:71:PHE:HA	2:H:85:GLN:O	2.14	0.47
2:F:4:GLU:N	2:F:4:GLU:OE1	2.48	0.47
2:D:33:SER:O	2:D:57:SER:HB3	2.15	0.46
3:E:150:LYS:HE2	3:E:155:LEU:HG	1.96	0.46
3:G:86:THR:OG1	3:G:104:LYS:HG2	2.15	0.46
2:D:103:ARG:O	2:D:108:ARG:HA	2.15	0.46
1:A:174:U:H2'	1:A:177:C:C6	2.50	0.46
3:E:31:SER:O	3:E:67:ARG:CZ	2.63	0.46
1:A:161:G:N7	2:H:106:ARG:NH1	2.62	0.46
3:E:33:ALA:O	3:E:91:GLN:HA	2.15	0.46
1:A:200:A:C6	1:B:200:A:C6	3.03	0.45
3:E:48:LEU:HA	3:E:59:VAL:HG21	1.98	0.45
2:H:86:MET:HB3	2:H:89:LEU:HD21	1.97	0.45
3:G:116:VAL:HA	3:G:136:LEU:O	2.17	0.45
3:L:30:VAL:HG11	3:L:91:GLN:CB	2.47	0.45
3:L:36:TRP:CH2	3:L:89:CYS:HB3	2.52	0.45
2:H:21:LEU:HD12	2:H:21:LEU:HA	1.84	0.44
2:D:4:GLU:CG	2:D:5:VAL:H	2.29	0.44
2:H:221:LYS:HE3	2:H:223:GLU:CD	2.38	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:143:ARG:HH21	3:L:164:VAL:HG11	1.81	0.43
3:L:25:ARG:NH1	3:L:71:ASP:OD1	2.51	0.43
2:F:79:LYS:O	2:F:81:THR:HG23	2.18	0.43
3:G:31:SER:O	3:G:67:ARG:CZ	2.66	0.43
3:E:109:ARG:HD2	3:E:172:SER:HB2	2.01	0.43
2:H:138:SER:OG	2:H:141:SER:O	2.37	0.43
1:A:237:C:H2'	1:A:238:C:C6	2.52	0.43
3:E:192:VAL:HG22	3:E:211:ASN:OD1	2.19	0.43
2:F:4:GLU:HG2	2:F:5:VAL:N	2.34	0.43
2:F:165:TRP:CH2	2:F:207:CYS:HB3	2.54	0.43
3:G:48:LEU:HA	3:G:59:VAL:HG21	2.00	0.43
2:D:79:LYS:O	2:D:81:THR:HG23	2.19	0.42
2:F:142:THR:HG1	3:G:117:PHE:HE2	1.60	0.42
1:B:155:A:H2'	1:B:156:C:C6	2.54	0.42
2:F:15:VAL:CG2	2:F:21:LEU:HD22	2.49	0.42
3:G:36:TRP:CD2	3:G:74:LEU:HB2	2.55	0.42
2:H:130:PRO:HB3	2:H:156:TYR:HB3	2.01	0.42
2:H:16:GLN:HA	2:H:123:SER:O	2.19	0.42
2:H:225:LYS:NZ	4:H:303:HOH:O	2.52	0.42
3:L:31:SER:O	3:L:67:ARG:CZ	2.67	0.42
2:H:109:ARG:HG2	3:L:50:TYR:CE2	2.54	0.42
3:E:91:GLN:OE1	3:E:93:SER:N	2.53	0.41
2:F:54:ILE:HA	2:F:60:SER:O	2.20	0.41
3:E:8:SER:OG	3:L:18:ASP:HB3	2.20	0.41
1:A:198:G:N2	1:A:201:A:OP2	2.47	0.41
3:E:195:CYS:O	3:E:207:THR:HA	2.21	0.41
1:A:237:C:H2'	1:A:238:C:H6	1.85	0.41
2:F:15:VAL:HG12	2:F:16:GLN:O	2.20	0.41
3:E:143:ARG:NH2	3:E:164:VAL:HG11	2.35	0.41
3:L:106:GLU:HG2	3:L:107:ILE:N	2.36	0.41
3:G:34:VAL:HA	3:G:90:GLN:O	2.22	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 X-ray entries followed by that with respect to entries of similar resolution.

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	
2	D	220/232 (95%)	213 (97%)	7 (3%)	0	100	100
2	F	220/232 (95%)	213 (97%)	7 (3%)	0	100	100
2	H	220/232 (95%)	212 (96%)	8 (4%)	0	100	100
3	E	213/215 (99%)	202 (95%)	11 (5%)	0	100	100
3	G	213/215 (99%)	202 (95%)	11 (5%)	0	100	100
3	L	213/215 (99%)	202 (95%)	11 (5%)	0	100	100
All	All	1299/1341 (97%)	1244 (96%)	55 (4%)	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 X-ray entries followed by that with respect to entries of similar resolution.

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	
2	D	185/195 (95%)	185 (100%)	0	100	100
2	F	185/195 (95%)	185 (100%)	0	100	100
2	H	185/195 (95%)	185 (100%)	0	100	100
3	E	190/190 (100%)	190 (100%)	0	100	100
3	G	190/190 (100%)	190 (100%)	0	100	100
3	L	190/190 (100%)	190 (100%)	0	100	100
All	All	1125/1155 (97%)	1125 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	67/68 (98%)	9 (13%)	0
1	B	67/68 (98%)	9 (13%)	0
1	C	67/68 (98%)	10 (14%)	0
All	All	201/204 (98%)	28 (13%)	0

All (28) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	154	A
1	A	158	G
1	A	161	G
1	A	164	U
1	A	233	G
1	A	234	U
1	A	235	G
1	A	236	C
1	A	248	C
1	B	154	A
1	B	158	G
1	B	161	G
1	B	164	U
1	B	233	G
1	B	234	U
1	B	235	G
1	B	236	C
1	B	248	C
1	C	154	A
1	C	158	G
1	C	161	G
1	C	164	U
1	C	219	C
1	C	233	G
1	C	234	U
1	C	235	G
1	C	236	C
1	C	248	C

There are no RNA pucker outliers to report.

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 [i](#)

There are no carbohydrates 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 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å²)	Q<0.9
1	A	68/68 (100%)	-0.11	0	100	100	62, 83, 130, 156	0
1	B	68/68 (100%)	-0.20	0	100	100	57, 83, 112, 136	0
1	C	68/68 (100%)	-0.18	0	100	100	70, 108, 198, 232	0
2	D	222/232 (95%)	0.69	8 (3%)	42	51	40, 59, 100, 175	0
2	F	222/232 (95%)	0.75	8 (3%)	42	51	42, 66, 107, 178	0
2	H	222/232 (95%)	0.68	6 (2%)	54	63	41, 58, 90, 150	0
3	E	215/215 (100%)	0.55	3 (1%)	75	82	37, 62, 99, 173	0
3	G	215/215 (100%)	0.61	9 (4%)	36	43	47, 77, 110, 147	0
3	L	215/215 (100%)	0.58	4 (1%)	66	75	41, 59, 91, 161	0
All	All	1515/1545 (98%)	0.54	38 (2%)	57	66	37, 65, 119, 232	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	142	THR	10.4
2	F	143	SER	6.5
2	D	142	THR	5.4
2	F	107	TYR	4.5
3	G	215	CYS	4.3
2	D	107	TYR	4.0
2	H	143	SER	3.9
2	H	138	SER	3.8
2	H	139	SER	3.7
3	G	30	VAL	3.7
2	F	144	GLY	3.6
3	G	20	VAL	3.5
2	D	139	SER	3.2
3	G	16	VAL	3.0
3	G	101	GLN	2.9

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Mol	Chain	Res	Type	RSRZ
2	D	144	GLY	2.8
3	L	155	LEU	2.6
2	D	140	LYS	2.5
3	G	3	ILE	2.5
2	F	139	SER	2.5
2	H	142	THR	2.4
2	F	145	GLY	2.4
3	E	214	GLU	2.3
3	G	76	ILE	2.3
2	H	151	CYS	2.2
2	D	143	SER	2.2
2	F	186	LEU	2.2
3	L	215	CYS	2.2
2	D	51	VAL	2.1
2	F	21	LEU	2.1
3	G	22	ILE	2.1
3	G	105	VAL	2.1
3	L	214	GLU	2.1
3	L	150	LYS	2.1
2	D	26	ALA	2.1
2	H	6	GLN	2.0
3	E	146	LYS	2.0
3	E	134	VAL	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.