



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:33 PM UTC

PDB ID : 7JRE / pdb_00007jre
Title : Crystal structure of EV-D68 2A protease C107A mutant
Authors : Liu, C.; Lee, M.-Y.; Liu, W.; Wang, J.
Deposited on : 2020-08-12
Resolution : 2.50 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 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:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

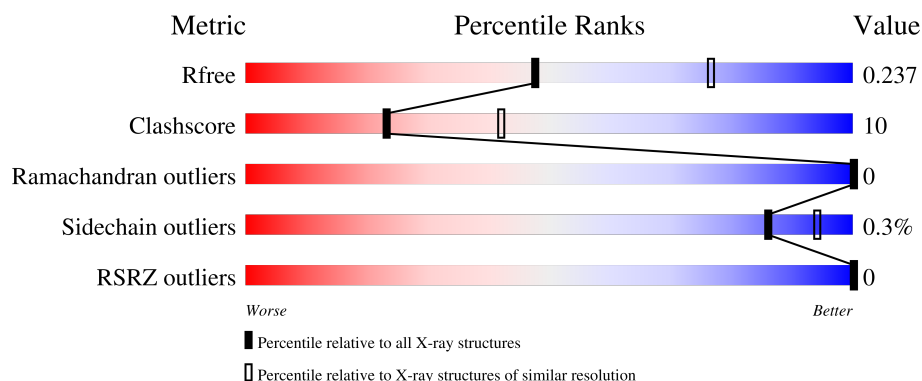
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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 of 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	147	 79% 15% 6%
1	B	147	 77% 16% 7%
1	C	147	 73% 21% 5%
1	D	147	 82% 12% • 5%
1	E	147	 80% 17% •

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Mol	Chain	Length	Quality of chain
1	F	147	78%15%7%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11815 atoms, of which 5600 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 protein called Protease 2A.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	138	Total 2017	C 668	H 976	N 181	O 187	S 5	0	0	0
1	B	137	Total 1916	C 643	H 916	N 174	O 178	S 5	0	0	0
1	C	139	Total 1914	C 652	H 894	N 179	O 184	S 5	0	0	0
1	D	139	Total 1954	C 656	H 931	N 174	O 188	S 5	0	0	0
1	E	142	Total 1983	C 658	H 947	N 183	O 190	S 5	0	0	0
1	F	137	Total 1966	C 659	H 936	N 181	O 185	S 5	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	107	ALA	CYS	engineered mutation	UNP A0A097BW19
B	107	ALA	CYS	engineered mutation	UNP A0A097BW19
C	107	ALA	CYS	engineered mutation	UNP A0A097BW19
D	107	ALA	CYS	engineered mutation	UNP A0A097BW19
E	107	ALA	CYS	engineered mutation	UNP A0A097BW19
F	107	ALA	CYS	engineered mutation	UNP A0A097BW19

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	1	Total 1	Zn 1	0	0
2	E	1	Total 1	Zn 1	0	0
2	F	1	Total 1	Zn 1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	18	Total 18	O 18	0	0
3	B	8	Total 8	O 8	0	0
3	C	13	Total 13	O 13	0	0
3	D	6	Total 6	O 6	0	0
3	E	8	Total 8	O 8	0	0
3	F	6	Total 6	O 6	0	0

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 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: Protease 2A

Chain A: 



• Molecule 1: Protease 2A

Chain B: 



• Molecule 1: Protease 2A

Chain C: 



• Molecule 1: Protease 2A

Chain D: 



• Molecule 1: Protease 2A

Chain E: 



• Molecule 1: Protease 2A

Chain F: 

GLY	PRO	GLY	F4	V7	F8	V9	G10	K13	V39	G46	R52	N56	T57	R63	W80	Q83	Y87	P88	A89	R90	L111	V117	F131	M136	L137	L138	W139	L140	ASP	THR	ASP	ALA	MET	GLU	GLN
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4 Data and refinement statistics

Property	Value	Source
Space group	P 4	Depositor
Cell constants a, b, c, α , β , γ	118.77Å 118.77Å 80.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.76 – 2.50 47.76 – 2.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (47.76-2.50) 99.9 (47.76-2.50)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.40 (at 2.51Å)	Xtriage
Refinement program	BUSTER, PHENIX dev_3893	Depositor
R, R_{free}	0.219 , 0.239 0.221 , 0.237	Depositor DCC
R_{free} test set	1979 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	62.8	Xtriage
Anisotropy	0.211	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 60.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.347 for h,-k,-l	Xtriage
Reported twinning fraction	0.370 for h,-k,-l	Depositor
Outliers	0 of 38883 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11815	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.15% 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](#)

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

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	0.28	0/1068	0.40	0/1454
1	B	0.34	0/1026	0.43	0/1400
1	C	0.26	0/1048	0.46	0/1428
1	D	0.47	0/1050	0.67	2/1434 (0.1%)
1	E	0.26	0/1062	0.39	0/1448
1	F	0.24	0/1058	0.39	0/1443
All	All	0.32	0/6312	0.47	2/8607 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	88	PRO	CA-N-CD	-10.95	96.67	112.00
1	D	89	ALA	CA-C-O	-6.30	114.65	120.95

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	1041	976	996	20	0
1	B	1000	916	929	22	0
1	C	1020	894	917	23	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1023	931	943	12	0
1	E	1036	947	966	21	0
1	F	1030	936	964	26	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	18	0	0	2	0
3	B	8	0	0	1	0
3	C	13	0	0	1	0
3	D	6	0	0	1	0
3	E	8	0	0	0	0
3	F	6	0	0	1	0
All	All	6215	5600	5715	122	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 10.

All (122) 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:83:GLN:HG2	1:F:88:PRO:HA	1.27	1.13
1:E:81:ILE:CG2	1:E:89:ALA:HB1	1.88	1.03
1:D:88:PRO:HD2	1:D:88:PRO:O	1.65	0.93
1:E:61:TYR:OH	1:E:66:ASP:OD1	1.91	0.89
1:C:35:SER:OG	1:C:94:ASN:OD1	1.94	0.86
1:E:81:ILE:HG22	1:E:89:ALA:HB1	1.59	0.83
1:D:61:TYR:O	3:D:301:HOH:O	1.97	0.82
1:B:14:ILE:HG23	1:B:39:VAL:HG22	1.66	0.78
1:F:83:GLN:HB2	1:F:89:ALA:HB2	1.64	0.78
1:F:83:GLN:CD	1:F:83:GLN:O	2.28	0.77
1:A:19:LEU:O	3:A:301:HOH:O	2.05	0.75
1:E:17:TYR:CD2	1:E:38:LEU:HD11	2.22	0.74
1:B:75:GLY:N	3:B:301:HOH:O	2.20	0.74
1:D:88:PRO:O	1:D:88:PRO:CD	2.37	0.72
1:E:52:ARG:CG	1:E:138:LEU:HD23	2.20	0.71
1:B:22:ILE:O	1:B:26:GLN:HG3	1.91	0.70
1:C:124:GLY:O	3:C:301:HOH:O	2.09	0.69
1:C:59:VAL:HB	1:D:59:VAL:HG23	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:17:TYR:HA	1:E:38:LEU:HD13	1.77	0.67
1:F:136:ASN:C	1:F:137:LEU:HD22	2.19	0.67
1:C:14:ILE:HG12	1:C:39:VAL:HG22	1.77	0.66
1:A:72:CYS:O	3:A:302:HOH:O	2.12	0.66
1:A:12:PHE:CE1	1:A:41:PRO:HG3	2.32	0.65
1:C:123:ALA:HB3	1:C:130:ALA:HB3	1.78	0.64
1:C:7:VAL:HG23	1:C:120:LEU:HD21	1.79	0.63
1:E:52:ARG:HG2	1:E:138:LEU:HD23	1.79	0.63
1:A:12:PHE:HE1	1:A:41:PRO:HG3	1.63	0.63
1:F:83:GLN:HG2	1:F:88:PRO:CA	2.18	0.63
1:A:90:ARG:HH11	1:A:90:ARG:HG3	1.64	0.62
1:A:47:ARG:HH11	1:A:47:ARG:HG3	1.63	0.62
1:A:36:ASP:OD2	1:A:122:THR:HG23	2.00	0.62
1:F:10:GLY:O	1:F:46:GLY:N	2.33	0.62
1:E:52:ARG:HG3	1:E:138:LEU:HD23	1.83	0.61
1:F:83:GLN:HB2	1:F:89:ALA:CB	2.32	0.60
1:C:128:ILE:HD12	1:C:128:ILE:C	2.26	0.59
1:F:7:VAL:HG12	1:F:9:VAL:HG23	1.84	0.59
1:A:90:ARG:HG3	1:A:90:ARG:NH1	2.18	0.59
1:D:10:GLY:O	1:D:46:GLY:N	2.37	0.57
1:D:9:VAL:HG21	1:D:50:ILE:HG13	1.86	0.57
1:E:81:ILE:HG23	1:E:89:ALA:HB1	1.79	0.56
1:B:122:THR:HG21	1:B:132:THR:OG1	2.04	0.56
1:F:83:GLN:HB2	1:F:89:ALA:H	1.71	0.56
1:C:132:THR:O	1:C:132:THR:HG23	2.06	0.55
1:F:80:TRP:CZ2	1:F:89:ALA:HB1	2.41	0.55
1:F:63:ARG:N	3:F:301:HOH:O	2.31	0.55
1:B:38:LEU:C	1:B:38:LEU:HD23	2.32	0.55
1:C:15:ILE:HD12	1:C:16:ASN:H	1.71	0.54
1:E:83:GLN:HG2	1:E:88:PRO:HA	1.89	0.54
1:F:7:VAL:HG21	1:F:117:VAL:HG11	1.89	0.54
1:D:125:GLY:O	1:D:128:ILE:HG13	2.07	0.54
1:B:24:GLU:HG3	1:B:40:THR:CB	2.38	0.53
1:F:7:VAL:CG1	1:F:9:VAL:HG23	2.38	0.53
1:A:29:ILE:HG23	1:A:41:PRO:HD3	1.89	0.53
1:C:128:ILE:HD12	1:C:128:ILE:O	2.10	0.52
1:F:83:GLN:HB2	1:F:89:ALA:N	2.24	0.51
1:B:3:GLY:O	1:B:61:TYR:CD2	2.64	0.51
1:B:50:ILE:HD12	1:B:134:ILE:HD12	1.92	0.51
1:A:46:GLY:O	1:A:47:ARG:HD3	2.11	0.51
1:F:7:VAL:CG2	1:F:117:VAL:HG11	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:38:LEU:HD23	1:B:39:VAL:N	2.25	0.51
1:B:25:ARG:HA	1:B:38:LEU:HD11	1.92	0.50
1:B:110:LEU:HD22	1:B:117:VAL:CG1	2.42	0.50
1:E:122:THR:OG1	1:E:132:THR:N	2.45	0.50
1:C:125:GLY:O	1:C:128:ILE:HG13	2.12	0.50
1:E:32:ASP:OD2	1:E:35:SER:OG	2.28	0.50
1:B:7:VAL:HG22	1:B:110:LEU:HD21	1.93	0.50
1:C:84:ASN:HB2	1:C:87:TYR:H	1.77	0.49
1:E:87:TYR:CE2	1:E:90:ARG:NH2	2.81	0.49
1:F:83:GLN:CB	1:F:89:ALA:N	2.75	0.49
1:C:13:LYS:HB2	1:C:42:ILE:HD13	1.94	0.49
1:F:87:TYR:CB	1:F:90:ARG:HH21	2.26	0.48
1:C:85:GLU:N	1:C:85:GLU:OE2	2.45	0.48
1:B:110:LEU:HD22	1:B:117:VAL:HG13	1.95	0.47
1:E:14:ILE:HG12	1:E:39:VAL:HG13	1.96	0.47
1:A:22:ILE:O	1:A:26:GLN:HG3	2.14	0.47
1:C:96:LEU:HB2	1:C:131:PHE:HB2	1.95	0.47
1:A:83:GLN:OE1	1:A:123:ALA:HB1	2.15	0.47
1:A:90:ARG:HH12	1:A:92:GLN:HG2	1.80	0.47
1:A:136:ASN:C	1:A:137:LEU:HD12	2.40	0.46
1:C:136:ASN:O	1:C:136:ASN:ND2	2.48	0.46
1:B:110:LEU:CD2	1:B:117:VAL:HG11	2.46	0.46
1:D:133:ASP:OD1	1:D:135:ARG:HG3	2.16	0.46
1:B:17:TYR:N	1:B:36:ASP:OD1	2.48	0.46
1:B:3:GLY:O	1:B:61:TYR:HD2	1.99	0.46
1:B:122:THR:CG2	1:B:132:THR:OG1	2.64	0.45
1:F:56:ASN:OD1	1:F:57:THR:HG23	2.17	0.45
1:A:52:ARG:NH2	1:A:138:LEU:O	2.50	0.45
1:C:32:ASP:HB3	1:C:37:VAL:HG12	1.99	0.45
1:F:111:LEU:HD22	1:F:131:PHE:CD2	2.52	0.44
1:E:17:TYR:HB3	1:E:87:TYR:OH	2.17	0.44
1:B:35:SER:O	1:B:132:THR:HG21	2.17	0.44
1:F:137:LEU:HD22	1:F:137:LEU:N	2.32	0.44
1:E:13:LYS:HE2	1:E:15:ILE:HD13	1.98	0.44
1:B:110:LEU:CD2	1:B:117:VAL:CG1	2.97	0.43
1:F:83:GLN:CG	1:F:88:PRO:HA	2.20	0.43
1:A:47:ARG:HG3	1:A:47:ARG:NH1	2.30	0.43
1:F:83:GLN:CB	1:F:89:ALA:H	2.30	0.43
1:A:90:ARG:NH1	1:A:92:GLN:HG2	2.33	0.43
1:B:18:HIS:HB3	1:B:86:TYR:CD2	2.54	0.43
1:F:83:GLN:HA	1:F:89:ALA:N	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:60:TYR:HA	1:D:110:LEU:O	2.19	0.42
1:E:81:ILE:HA	1:E:90:ARG:O	2.18	0.42
1:C:121:LEU:O	1:C:130:ALA:O	2.37	0.42
1:A:17:TYR:HB3	1:A:36:ASP:HA	2.01	0.42
1:D:111:LEU:HD13	1:D:131:PHE:CE2	2.55	0.42
1:E:113:CYS:HB3	1:E:118:ILE:HD11	2.01	0.42
1:D:56:ASN:OD1	1:D:56:ASN:C	2.63	0.42
1:E:83:GLN:OE1	1:E:84:ASN:N	2.53	0.41
1:C:73:PHE:CD2	1:C:96:LEU:HD22	2.55	0.41
1:C:122:THR:O	1:C:123:ALA:HB2	2.20	0.41
1:A:87:TYR:HB3	1:A:90:ARG:HE	1.86	0.41
1:B:32:ASP:OD2	1:B:35:SER:OG	2.26	0.41
1:C:87:TYR:CE2	1:C:90:ARG:NH2	2.89	0.41
1:F:52:ARG:NH2	1:F:138:LEU:O	2.53	0.41
1:E:88:PRO:O	1:E:90:ARG:HD3	2.21	0.41
1:A:82:GLU:O	1:A:82:GLU:HG2	2.21	0.41
1:F:83:GLN:CB	1:F:89:ALA:HB2	2.44	0.41
1:C:59:VAL:HG11	1:D:113:CYS:HA	2.02	0.40
1:E:79:GLN:CD	1:E:80:TRP:H	2.29	0.40
1:B:35:SER:O	1:B:132:THR:CG2	2.69	0.40
1:C:4:PHE:O	1:C:4:PHE:CG	2.74	0.40
1:F:13:LYS:O	1:F:39:VAL:HA	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 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	
1	A	136/147 (92%)	129 (95%)	7 (5%)	0	100	100
1	B	135/147 (92%)	129 (96%)	6 (4%)	0	100	100
1	C	137/147 (93%)	127 (93%)	10 (7%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	137/147 (93%)	133 (97%)	4 (3%)	0	100	100
1	E	140/147 (95%)	136 (97%)	4 (3%)	0	100	100
1	F	135/147 (92%)	123 (91%)	12 (9%)	0	100	100
All	All	820/882 (93%)	777 (95%)	43 (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 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	
1	A	102/114 (90%)	102 (100%)	0	100	100
1	B	90/114 (79%)	90 (100%)	0	100	100
1	C	92/114 (81%)	91 (99%)	1 (1%)	65	84
1	D	96/114 (84%)	95 (99%)	1 (1%)	68	86
1	E	97/114 (85%)	97 (100%)	0	100	100
1	F	98/114 (86%)	98 (100%)	0	100	100
All	All	575/684 (84%)	573 (100%)	2 (0%)	86	94

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	38	LEU
1	D	47	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	92	GLN
1	B	16	ASN
1	B	26	GLN

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Mol	Chain	Res	Type
1	C	136	ASN
1	D	16	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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(Å ²)	Q<0.9
1	A	138/147 (93%)	-0.73	0 100 100	43, 60, 76, 82	0
1	B	137/147 (93%)	-0.60	0 100 100	38, 68, 86, 95	0
1	C	139/147 (94%)	-0.51	0 100 100	60, 76, 90, 93	0
1	D	139/147 (94%)	-0.52	0 100 100	56, 79, 92, 99	0
1	E	142/147 (96%)	-0.60	0 100 100	46, 65, 83, 90	0
1	F	137/147 (93%)	-0.60	0 100 100	55, 72, 91, 103	0
All	All	832/882 (94%)	-0.59	0 100 100	38, 71, 89, 103	0

There are no RSRZ outliers to report.

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 oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ZN	A	201	1/1	1.00	0.02	54,54,54,54	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ZN	B	201	1/1	1.00	0.01	54,54,54,54	0
2	ZN	C	201	1/1	1.00	0.01	74,74,74,74	0
2	ZN	D	201	1/1	1.00	0.02	64,64,64,64	0
2	ZN	E	201	1/1	1.00	0.01	59,59,59,59	0
2	ZN	F	201	1/1	1.00	0.02	58,58,58,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.