



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 10, 2023 – 01:56 AM EDT

PDB ID : 7MG0
Title : Crystal structure of EV-D68 2A protease
Authors : Liu, C.; Lee, M.-Y.; Liu, W.; Wang, J.
Deposited on : 2021-04-12
Resolution : 3.50 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.35.1
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) : 2.35.1

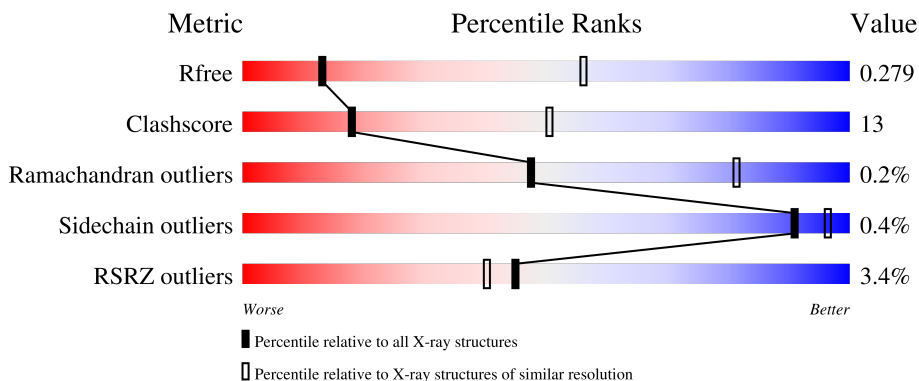
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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.



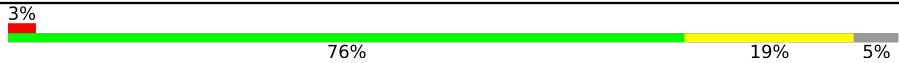
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1659 (3.60-3.40)
Clashscore	141614	1036 (3.58-3.42)
Ramachandran outliers	138981	1005 (3.58-3.42)
Sidechain outliers	138945	1006 (3.58-3.42)
RSRZ outliers	127900	1559 (3.60-3.40)

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	<div style="display: flex; align-items: center;"> <div style="width: 4%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 78%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 18%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div> <p style="margin-left: 40px;">78% 18% . .</p>
1	B	147	<div style="display: flex; align-items: center;"> <div style="width: 75%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 20%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: grey;"></div> </div> <p style="margin-left: 40px;">75% 20% . .</p>
1	C	147	<div style="display: flex; align-items: center;"> <div style="width: 7%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 69%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 24%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: grey;"></div> </div> <p style="margin-left: 40px;">69% 24% . 5%</p>
1	D	147	<div style="display: flex; align-items: center;"> <div style="width: 5%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 73%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 21%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 6%; height: 10px; background-color: grey;"></div> </div> <p style="margin-left: 40px;">73% 21% 6%</p>
1	E	147	<div style="display: flex; align-items: center;"> <div style="width: 0%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 76%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 20%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 4%; height: 10px; background-color: grey;"></div> </div> <p style="margin-left: 40px;">76% 20% .</p>

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Mol	Chain	Length	Quality of chain
1	F	147	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into four segments: a small red segment on the left labeled '3%', a large green segment labeled '76%', a yellow segment labeled '19%', and a small grey segment on the right labeled '5%'.</p>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11167 atoms, of which 5169 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
			Total	C	H	N	O	S			
1	E	142	1933	657	899	183	188	6	0	0	0
1	A	141	1813	621	828	175	183	6	0	0	0
1	B	141	1892	650	872	177	187	6	0	0	0
1	C	139	1840	618	863	175	178	6	0	0	0
1	D	138	1835	620	856	175	178	6	0	0	0
1	F	140	1836	623	851	173	183	6	0	0	0

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

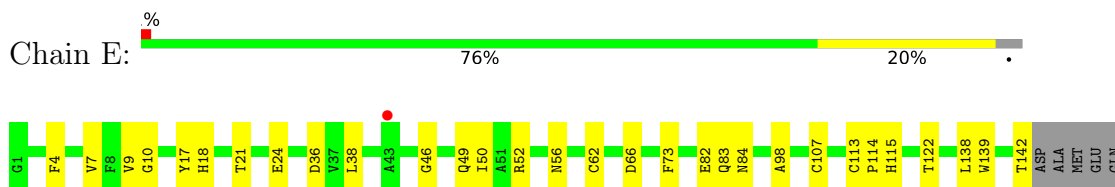
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	E	2	Total O 2 2	0	0
3	A	1	Total O 1 1	0	0
3	B	2	Total O 2 2	0	0
3	C	1	Total O 1 1	0	0
3	D	2	Total O 2 2	0	0
3	F	4	Total O 4 4	0	0

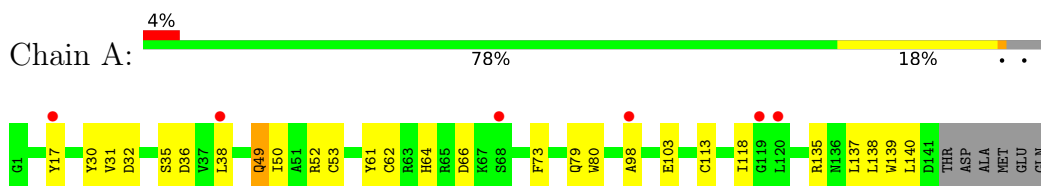
3 Residue-property plots [i](#)

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.

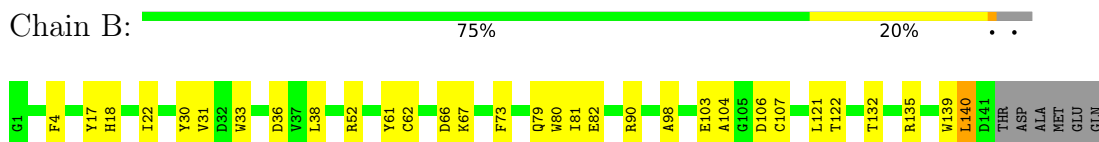
- Molecule 1: Protease 2A



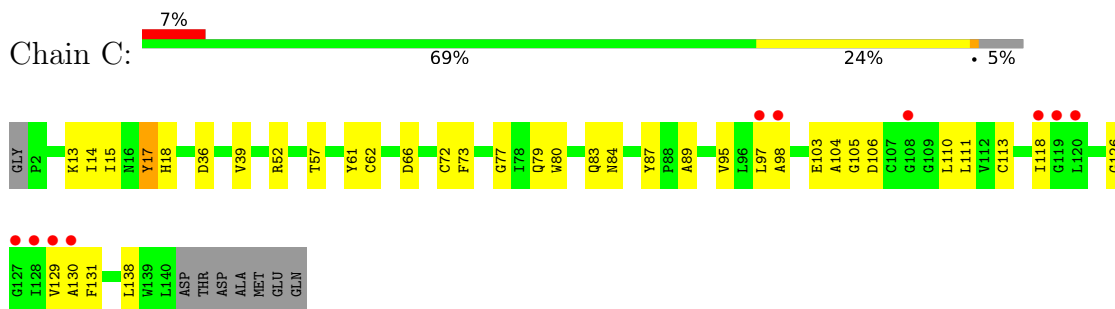
- Molecule 1: Protease 2A



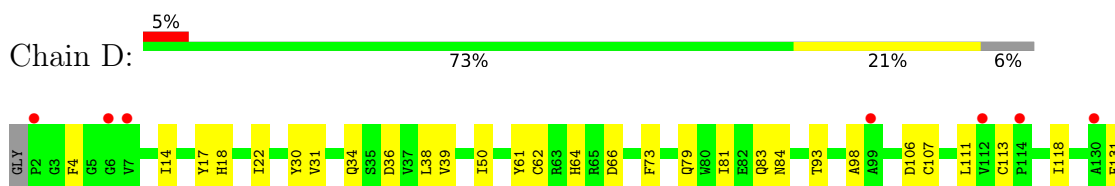
- Molecule 1: Protease 2A

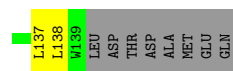


- Molecule 1: Protease 2A



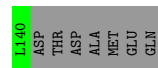
- Molecule 1: Protease 2A





- Molecule 1: Protease 2A

Chain F: 3% 76% 19% 5%



4 Data and refinement statistics i

Property	Value	Source
Space group	P 4	Depositor
Cell constants a, b, c, α , β , γ	118.16Å 118.16Å 81.73Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.88 – 3.50 47.88 – 3.20	Depositor EDS
% Data completeness (in resolution range)	100.0 (47.88-3.50) 96.8 (47.88-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.28 (at 3.19Å)	Xtrriage
Refinement program	PHENIX dev_4189	Depositor
R, R_{free}	0.238 , 0.285 0.235 , 0.279	Depositor DCC
R_{free} test set	1883 reflections (10.04%)	wwPDB-VP
Wilson B-factor (Å ²)	107.0	Xtrriage
Anisotropy	0.080	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 104.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtrriage
Estimated twinning fraction	0.468 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	11167	wwPDB-VP
Average B, all atoms (Å ²)	128.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.14% 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	0.27	0/1005	0.56	0/1371
1	B	0.26	0/1046	0.54	0/1427
1	C	0.28	0/1000	0.58	0/1365
1	D	0.26	0/1002	0.51	0/1366
1	E	0.26	0/1060	0.53	0/1445
1	F	0.30	0/1007	0.56	0/1375
All	All	0.27	0/6120	0.55	0/8349

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	985	828	912	25	0
1	B	1020	872	944	33	0
1	C	977	863	893	27	0
1	D	979	856	902	31	0
1	E	1034	899	964	24	0
1	F	985	851	909	22	0
2	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
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	1	0	0	0	0
3	B	2	0	0	0	0
3	C	1	0	0	0	0
3	D	2	0	0	3	0
3	E	2	0	0	1	0
3	F	4	0	0	0	0
All	All	5998	5169	5524	155	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 (155) 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:B:31:VAL:HG22	1:F:31:VAL:HG22	1.51	0.91
1:A:113:CYS:HB3	1:A:118:ILE:HD11	1.50	0.90
1:B:61:TYR:OH	1:B:66:ASP:OD1	1.98	0.81
1:D:31:VAL:HG12	1:D:38:LEU:HG	1.63	0.79
1:E:82:GLU:OE1	1:E:84:ASN:ND2	2.17	0.76
1:A:31:VAL:HG22	1:D:31:VAL:HG22	1.68	0.74
1:C:113:CYS:HB3	1:C:118:ILE:HD11	1.72	0.72
1:A:61:TYR:OH	1:A:66:ASP:OD1	2.08	0.71
1:A:32:ASP:OD2	1:A:35:SER:OG	2.05	0.71
1:B:139:TRP:O	1:B:140:LEU:HB2	1.90	0.70
1:D:138:LEU:HA	3:D:301:HOH:O	1.90	0.69
1:A:73:PHE:HB3	1:A:98:ALA:HB2	1.75	0.69
1:D:83:GLN:NE2	1:D:84:ASN:O	2.29	0.66
1:D:111:LEU:HD22	1:D:131:PHE:HD2	1.61	0.66
1:C:113:CYS:HB3	1:C:118:ILE:CD1	2.26	0.65
1:E:139:TRP:NE1	1:E:142:THR:HA	2.12	0.65
1:D:79:GLN:HA	1:D:79:GLN:OE1	1.97	0.64
1:A:30:TYR:HA	1:D:31:VAL:HG23	1.79	0.64
1:E:73:PHE:HB3	1:E:98:ALA:HB2	1.78	0.64
1:F:31:VAL:HG12	1:F:38:LEU:HG	1.80	0.63
1:B:30:TYR:HA	1:F:31:VAL:HG23	1.81	0.62
1:C:103:GLU:HG3	1:C:104:ALA:N	2.15	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:73:PHE:HB3	1:D:98:ALA:HB2	1.83	0.60
1:F:122:THR:HG1	1:F:132:THR:HG1	1.47	0.60
1:D:64:HIS:ND1	1:D:106:ASP:OD2	2.35	0.60
1:A:53:CYS:O	1:A:135:ARG:NH1	2.34	0.60
1:E:9:VAL:HG21	1:E:50:ILE:HD11	1.83	0.60
1:A:113:CYS:CB	1:A:118:ILE:HD11	2.27	0.59
1:D:113:CYS:HB3	1:D:118:ILE:HD11	1.84	0.59
1:F:81:ILE:HG23	1:F:81:ILE:O	2.02	0.59
1:E:83:GLN:NE2	1:E:84:ASN:O	2.36	0.58
1:C:77:GLY:O	1:C:95:VAL:HB	2.02	0.58
1:C:111:LEU:HD22	1:C:131:PHE:CD2	2.39	0.58
1:D:50:ILE:HD11	1:D:138:LEU:HD12	1.87	0.57
1:D:137:LEU:C	3:D:301:HOH:O	2.42	0.57
1:A:137:LEU:O	1:A:138:LEU:HB2	2.05	0.56
1:D:81:ILE:O	1:D:81:ILE:HG23	2.05	0.56
1:A:62:CYS:O	1:A:66:ASP:N	2.38	0.56
1:C:73:PHE:HB3	1:C:98:ALA:HB2	1.87	0.56
1:F:113:CYS:HB3	1:F:118:ILE:HD11	1.87	0.56
1:C:83:GLN:HB2	1:C:89:ALA:HB2	1.86	0.56
1:B:73:PHE:HB3	1:B:98:ALA:HB2	1.87	0.56
1:C:111:LEU:HD22	1:C:131:PHE:HD2	1.71	0.56
1:E:56:ASN:OD1	1:B:67:LYS:NZ	2.40	0.55
1:C:103:GLU:HG3	1:C:105:GLY:H	1.71	0.55
1:A:31:VAL:HG23	1:D:30:TYR:HA	1.89	0.54
1:B:18:HIS:CE1	1:B:107:CYS:HB3	2.41	0.54
1:B:81:ILE:HG23	1:B:90:ARG:N	2.22	0.54
1:D:14:ILE:HG12	1:D:39:VAL:HG22	1.89	0.54
1:D:73:PHE:CE2	1:D:111:LEU:HD21	2.42	0.54
1:A:52:ARG:NH2	1:A:137:LEU:O	2.41	0.54
1:B:80:TRP:O	1:B:82:GLU:OE1	2.26	0.54
1:F:20:ALA:CB	1:F:38:LEU:HD22	2.38	0.53
1:D:138:LEU:CA	3:D:301:HOH:O	2.54	0.53
1:C:14:ILE:HG12	1:C:39:VAL:HG22	1.91	0.53
1:F:73:PHE:HB3	1:F:98:ALA:HB2	1.90	0.53
1:D:17:TYR:N	1:D:36:ASP:OD1	2.41	0.52
1:A:17:TYR:N	1:A:36:ASP:OD1	2.42	0.52
1:C:79:GLN:CG	1:C:80:TRP:N	2.72	0.52
1:B:62:CYS:O	1:B:66:ASP:N	2.42	0.52
1:C:84:ASN:OD1	1:C:87:TYR:N	2.33	0.52
1:E:62:CYS:O	1:E:66:ASP:N	2.41	0.52
1:B:18:HIS:HE1	1:B:107:CYS:HB3	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:52:ARG:HG2	1:C:138:LEU:HD13	1.93	0.51
1:E:4:PHE:O	1:E:4:PHE:CD2	2.64	0.51
1:D:4:PHE:HB3	1:D:61:TYR:HD2	1.76	0.50
1:B:122:THR:OG1	1:B:132:THR:OG1	2.16	0.50
1:E:49:GLN:NE2	1:E:115:HIS:O	2.44	0.50
1:C:110:LEU:HD13	1:C:111:LEU:N	2.27	0.49
1:B:52:ARG:HA	1:B:140:LEU:CD2	2.42	0.49
1:B:106:ASP:O	1:B:121:LEU:HD23	2.12	0.49
1:B:31:VAL:HG23	1:F:30:TYR:HA	1.94	0.49
1:A:64:HIS:CD2	1:A:103:GLU:HG3	2.48	0.48
1:A:79:GLN:HG3	1:A:80:TRP:H	1.77	0.48
1:B:52:ARG:HA	1:B:140:LEU:HD23	1.95	0.48
1:C:103:GLU:N	1:C:106:ASP:OD2	2.47	0.48
1:A:103:GLU:OE1	1:A:103:GLU:HA	2.14	0.48
1:E:21:THR:HG22	1:E:24:GLU:OE1	2.13	0.48
1:D:64:HIS:N	1:D:106:ASP:OD1	2.41	0.48
1:E:52:ARG:NH2	1:E:138:LEU:O	2.46	0.48
1:E:36:ASP:OD2	1:E:122:THR:HG23	2.14	0.47
1:B:4:PHE:HD1	1:B:4:PHE:O	1.97	0.47
1:E:17:TYR:HA	1:E:38:LEU:HD13	1.95	0.47
1:B:4:PHE:HD2	1:B:61:TYR:HE2	1.63	0.47
1:E:18:HIS:CE1	1:E:107:CYS:HB3	2.50	0.47
1:E:17:TYR:N	1:E:36:ASP:OD1	2.48	0.47
1:B:139:TRP:O	1:B:140:LEU:CB	2.62	0.47
1:D:4:PHE:HB3	1:D:61:TYR:CD2	2.50	0.47
1:D:62:CYS:O	1:D:66:ASP:N	2.48	0.46
1:D:4:PHE:O	1:D:4:PHE:HD1	1.99	0.46
1:D:111:LEU:HD22	1:D:131:PHE:CD2	2.45	0.46
1:F:51:ALA:O	1:F:135:ARG:HG2	2.16	0.46
1:B:17:TYR:N	1:B:36:ASP:OD1	2.46	0.46
1:F:17:TYR:HA	1:F:38:LEU:HD13	1.98	0.46
1:E:7:VAL:HG12	1:E:9:VAL:HG23	1.99	0.45
1:F:138:LEU:H	1:F:138:LEU:HD12	1.80	0.45
1:F:20:ALA:HB1	1:F:38:LEU:HD22	1.98	0.45
1:A:62:CYS:SG	1:A:64:HIS:ND1	2.90	0.45
1:E:21:THR:HA	3:E:301:HOH:O	2.17	0.44
1:A:139:TRP:O	1:A:140:LEU:HB2	2.17	0.44
1:C:113:CYS:CB	1:C:118:ILE:HD11	2.43	0.44
1:E:73:PHE:CB	1:E:98:ALA:HB2	2.47	0.44
1:A:64:HIS:CE1	1:A:103:GLU:HG2	2.52	0.44
1:D:22:ILE:H	1:D:22:ILE:HD12	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:79:GLN:OE1	1:C:97:LEU:HD22	2.18	0.44
1:E:139:TRP:CE3	1:E:139:TRP:HA	2.52	0.44
1:B:79:GLN:HG3	1:B:80:TRP:H	1.81	0.44
1:A:31:VAL:HG12	1:A:38:LEU:HG	2.00	0.44
1:D:34:GLN:HB3	1:D:93:THR:HG22	1.99	0.44
1:E:17:TYR:CD2	1:E:38:LEU:HD11	2.52	0.44
1:A:137:LEU:HD23	1:A:137:LEU:HA	1.90	0.44
1:B:139:TRP:O	1:B:140:LEU:HD13	2.17	0.44
1:C:62:CYS:O	1:C:66:ASP:N	2.51	0.44
1:B:31:VAL:HG12	1:B:38:LEU:HG	2.00	0.43
1:D:4:PHE:O	1:D:4:PHE:CD1	2.71	0.43
1:E:10:GLY:O	1:E:46:GLY:N	2.48	0.43
1:E:113:CYS:HB2	1:E:114:PRO:HD2	2.00	0.43
1:B:79:GLN:HG3	1:B:80:TRP:N	2.33	0.43
1:B:79:GLN:CG	1:B:80:TRP:H	2.31	0.43
1:E:9:VAL:HG21	1:E:50:ILE:CD1	2.49	0.43
1:A:62:CYS:SG	1:A:64:HIS:CE1	3.12	0.43
1:B:4:PHE:O	1:B:4:PHE:CD1	2.72	0.43
1:B:135:ARG:O	1:B:140:LEU:CD1	2.67	0.43
1:D:113:CYS:HB3	1:D:118:ILE:CD1	2.47	0.43
1:B:22:ILE:HD12	1:B:22:ILE:H	1.84	0.43
1:C:17:TYR:N	1:C:36:ASP:OD1	2.52	0.43
1:A:64:HIS:CG	1:A:103:GLU:HG3	2.54	0.42
1:B:103:GLU:HG3	1:B:104:ALA:N	2.33	0.42
1:C:129:VAL:HG22	1:C:130:ALA:N	2.34	0.42
1:B:17:TYR:HA	1:B:38:LEU:CD1	2.49	0.42
1:F:24:GLU:HG2	1:F:40:THR:HG21	2.00	0.42
1:C:13:LYS:NZ	1:C:15:ILE:HG21	2.35	0.42
1:A:17:TYR:HD1	1:A:38:LEU:HD11	1.85	0.42
1:C:57:THR:HG22	1:C:72:CYS:HB2	2.02	0.42
1:F:110:LEU:HD23	1:F:120:LEU:CD2	2.50	0.42
1:B:17:TYR:CD1	1:B:38:LEU:HD11	2.55	0.42
1:D:62:CYS:SG	1:D:64:HIS:CE1	3.13	0.42
1:C:18:HIS:O	1:C:18:HIS:HD2	2.03	0.41
1:F:32:ASP:OD2	1:F:35:SER:OG	2.15	0.41
1:F:64:HIS:ND1	1:F:106:ASP:OD2	2.53	0.41
1:A:49:GLN:HG3	1:A:50:ILE:N	2.36	0.41
1:F:113:CYS:HB3	1:F:118:ILE:CD1	2.50	0.41
1:D:18:HIS:CE1	1:D:107:CYS:HB3	2.55	0.41
1:F:32:ASP:O	1:F:36:ASP:N	2.54	0.41
1:C:83:GLN:HG3	1:C:89:ALA:N	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:7:VAL:CG1	1:E:9:VAL:HG23	2.51	0.41
1:C:61:TYR:OH	1:C:66:ASP:OD1	2.24	0.41
1:C:79:GLN:HG2	1:C:80:TRP:N	2.36	0.41
1:B:4:PHE:CD2	1:B:61:TYR:HE2	2.39	0.40
1:B:33:TRP:O	1:B:90:ARG:NH2	2.54	0.40
1:F:103:GLU:HG3	1:F:104:ALA:H	1.86	0.40
1:D:17:TYR:HA	1:D:38:LEU:CD1	2.51	0.40
1:F:100:GLY:N	1:F:127:GLY:O	2.48	0.40
1:C:79:GLN:OE1	1:C:97:LEU:HD13	2.22	0.40
1:F:17:TYR:N	1:F:36:ASP:OD1	2.52	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	139/147 (95%)	127 (91%)	12 (9%)	0	100	100
1	B	139/147 (95%)	131 (94%)	7 (5%)	1 (1%)	22	61
1	C	137/147 (93%)	120 (88%)	16 (12%)	1 (1%)	22	61
1	D	136/147 (92%)	127 (93%)	9 (7%)	0	100	100
1	E	140/147 (95%)	132 (94%)	8 (6%)	0	100	100
1	F	138/147 (94%)	130 (94%)	8 (6%)	0	100	100
All	All	829/882 (94%)	767 (92%)	60 (7%)	2 (0%)	47	81

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	140	LEU
1	C	126	GLY

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 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	91/115 (79%)	90 (99%)	1 (1%)	73	88
1	B	97/115 (84%)	97 (100%)	0	100	100
1	C	89/115 (77%)	88 (99%)	1 (1%)	73	88
1	D	91/115 (79%)	91 (100%)	0	100	100
1	E	98/115 (85%)	98 (100%)	0	100	100
1	F	92/115 (80%)	92 (100%)	0	100	100
All	All	558/690 (81%)	556 (100%)	2 (0%)	91	96

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

Mol	Chain	Res	Type
1	A	49	GLN
1	C	17	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	84	ASN
1	B	18	HIS
1	C	18	HIS

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

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 [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(Å ²)	Q<0.9
1	A	141/147 (95%)	0.16	6 (4%) 35 31	77, 119, 150, 166	0
1	B	141/147 (95%)	0.11	0 100 100	67, 100, 143, 174	0
1	C	139/147 (94%)	0.47	10 (7%) 15 15	74, 131, 168, 212	0
1	D	138/147 (93%)	0.31	7 (5%) 28 25	77, 129, 161, 187	0
1	E	142/147 (96%)	0.18	1 (0%) 87 83	67, 98, 145, 178	0
1	F	140/147 (95%)	0.20	5 (3%) 42 38	78, 121, 151, 178	0
All	All	841/882 (95%)	0.24	29 (3%) 45 40	67, 117, 157, 212	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	89	ALA	5.1
1	A	119	GLY	4.1
1	C	129	VAL	3.5
1	D	6	GLY	3.3
1	C	98	ALA	3.2
1	D	114	PRO	3.2
1	F	110	LEU	3.2
1	C	128	ILE	3.2
1	E	43	ALA	3.1
1	F	3	GLY	3.0
1	D	2	PRO	2.9
1	A	38	LEU	2.9
1	A	17	TYR	2.9
1	C	127	GLY	2.8
1	A	120	LEU	2.8
1	C	119	GLY	2.7
1	D	99	ALA	2.6
1	C	130	ALA	2.5
1	C	108	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	F	4	PHE	2.4
1	A	68	SER	2.2
1	C	120	LEU	2.2
1	F	111	LEU	2.2
1	A	98	ALA	2.2
1	D	130	ALA	2.2
1	D	112	VAL	2.2
1	D	7	VAL	2.1
1	C	97	LEU	2.0
1	C	118	ILE	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 monosaccharides 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	0.93	0.24	148,148,148,148	0
2	ZN	C	201	1/1	0.94	0.27	138,138,138,138	0
2	ZN	F	201	1/1	0.95	0.18	122,122,122,122	0
2	ZN	E	201	1/1	0.96	0.30	99,99,99,99	0
2	ZN	D	201	1/1	0.99	0.24	99,99,99,99	0
2	ZN	B	201	1/1	0.99	0.27	58,58,58,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.