



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:20 PM UTC

PDB ID : 3B48 / pdb_00003b48
Title : Crystal structure of unknown function protein EF1359
Authors : Chang, C.; Li, H.; Moy, S.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2007-10-23
Resolution : 2.21 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

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Mol	Chain	Length	Quality of chain
1	F	135	3% 76% 22%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6128 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 protein called Uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	129	Total	C	N	O	Se	0	0	0
			963	599	156	202	6			
1	B	132	Total	C	N	O	Se	0	1	0
			991	616	160	209	6			
1	C	134	Total	C	N	O	Se	0	0	0
			999	621	161	211	6			
1	D	130	Total	C	N	O	Se	0	0	0
			961	600	154	201	6			
1	E	130	Total	C	N	O	Se	0	0	0
			966	603	155	202	6			
1	F	135	Total	C	N	O	Se	3	1	0
			1012	627	164	215	6			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP Q835L8
A	-1	ASN	-	expression tag	UNP Q835L8
A	0	ALA	-	expression tag	UNP Q835L8
B	-2	SER	-	expression tag	UNP Q835L8
B	-1	ASN	-	expression tag	UNP Q835L8
B	0	ALA	-	expression tag	UNP Q835L8
C	-2	SER	-	expression tag	UNP Q835L8
C	-1	ASN	-	expression tag	UNP Q835L8
C	0	ALA	-	expression tag	UNP Q835L8
D	-2	SER	-	expression tag	UNP Q835L8
D	-1	ASN	-	expression tag	UNP Q835L8
D	0	ALA	-	expression tag	UNP Q835L8
E	-2	SER	-	expression tag	UNP Q835L8
E	-1	ASN	-	expression tag	UNP Q835L8
E	0	ALA	-	expression tag	UNP Q835L8
F	-2	SER	-	expression tag	UNP Q835L8
F	-1	ASN	-	expression tag	UNP Q835L8

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Chain	Residue	Modelled	Actual	Comment	Reference
F	0	ALA	-	expression tag	UNP Q835L8

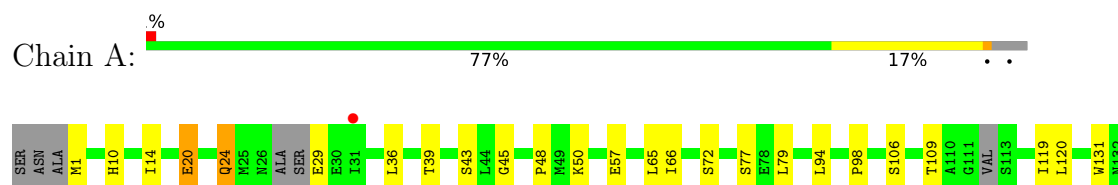
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	44	Total O 44 44	0	0
2	B	47	Total O 47 47	0	0
2	C	45	Total O 45 45	0	0
2	D	30	Total O 30 30	0	0
2	E	33	Total O 33 33	0	0
2	F	37	Total O 37 37	0	0

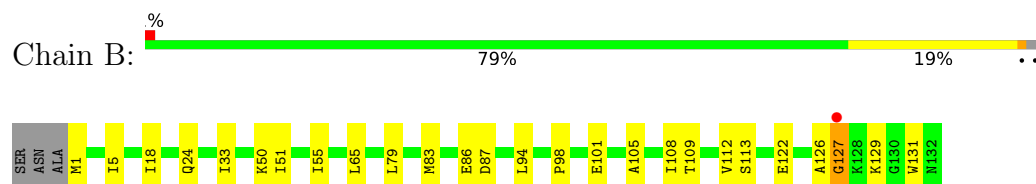
3 Residue-property plots [i](#)

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.

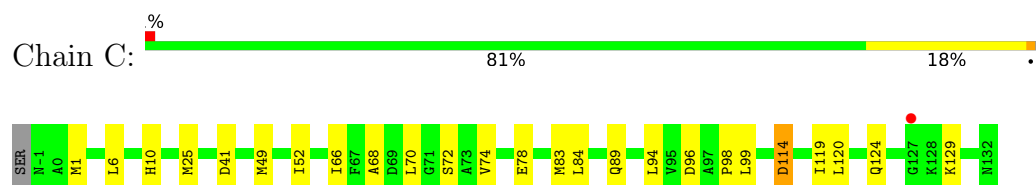
- Molecule 1: Uncharacterized protein



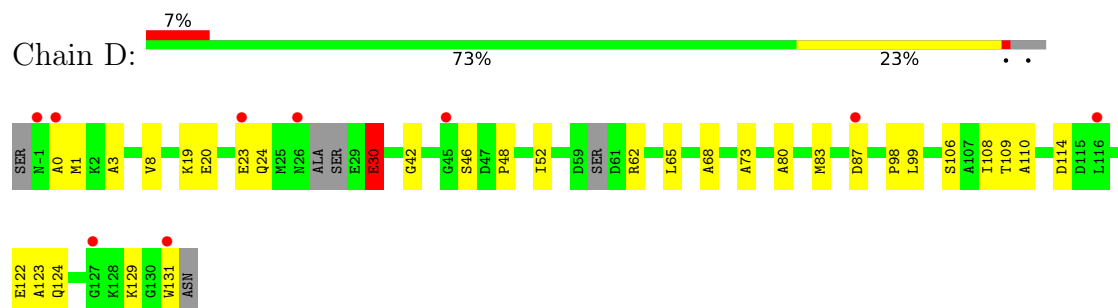
- Molecule 1: Uncharacterized protein



- Molecule 1: Uncharacterized protein

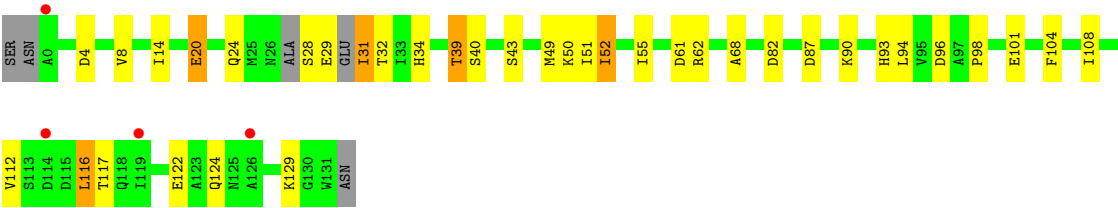


- Molecule 1: Uncharacterized protein

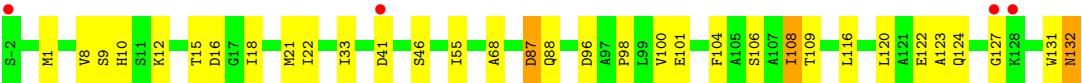
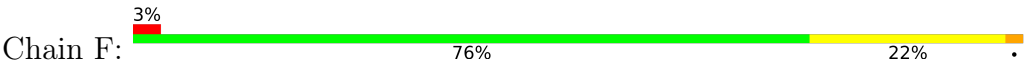


- Molecule 1: Uncharacterized protein





● Molecule 1: Uncharacterized protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	60.08Å 102.11Å 60.45Å 90.00° 96.57° 90.00°	Depositor
Resolution (Å)	44.99 – 2.21 44.99 – 2.21	Depositor EDS
% Data completeness (in resolution range)	94.5 (44.99-2.21) 94.5 (44.99-2.21)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.15 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.201 , 0.246 0.200 , 0.242	Depositor DCC
R_{free} test set	1713 reflections (4.71%)	wwPDB-VP
Wilson B-factor (Å ²)	30.3	Xtriage
Anisotropy	0.566	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 39.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.042 for l,-k,h	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6128	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.74% 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

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	0.89	0/968	1.05	0/1294
1	B	0.89	0/998	1.07	3/1337 (0.2%)
1	C	0.88	0/1005	1.13	3/1345 (0.2%)
1	D	0.78	0/965	1.03	2/1292 (0.2%)
1	E	0.84	0/970	1.03	0/1297
1	F	0.83	0/1018	1.08	1/1363 (0.1%)
All	All	0.85	0/5924	1.07	9/7928 (0.1%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	126	ALA	N-CA-C	10.13	124.15	111.69
1	D	30	GLU	N-CA-C	-7.42	103.80	114.12
1	B	127	GLY	N-CA-C	7.33	123.37	113.99
1	C	70	LEU	N-CA-C	5.77	117.55	108.96
1	B	113	SER	N-CA-C	5.56	118.53	109.24
1	D	99	LEU	N-CA-C	5.44	117.66	111.02
1	F	108	ILE	N-CA-C	5.12	115.34	110.42
1	C	99	LEU	N-CA-C	5.03	117.16	111.02
1	C	114	ASP	N-CA-C	5.01	118.43	112.72

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	963	0	928	22	0
1	B	991	0	962	12	0
1	C	999	0	969	26	0
1	D	961	0	920	28	0
1	E	966	0	934	27	0
1	F	1012	0	977	26	0
2	A	44	0	0	1	0
2	B	47	0	0	1	0
2	C	45	0	0	1	0
2	D	30	0	0	1	0
2	E	33	0	0	5	0
2	F	37	0	0	2	0
All	All	6128	0	5690	119	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 (119) 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:D:1:MSE:HE3	1:D:3:ALA:CB	1.67	1.23
1:A:79:LEU:CD1	1:C:49:MSE:HE3	1.67	1.22
1:D:80:ALA:HA	1:D:83:MSE:HE2	1.30	1.11
1:D:1:MSE:HE3	1:D:3:ALA:HB2	1.13	1.08
1:A:79:LEU:HD11	1:C:49:MSE:HE3	1.37	1.02
1:C:49:MSE:CE	1:C:52:ILE:HG13	1.98	0.93
1:E:29:GLU:HG2	1:E:31:ILE:HD13	1.50	0.90
1:C:49:MSE:HE2	1:C:52:ILE:HG13	1.54	0.90
1:D:52:ILE:CG1	1:D:83:MSE:HE3	2.05	0.87
1:E:29:GLU:CG	1:E:31:ILE:HD13	2.06	0.85
1:E:124:GLN:HG2	2:E:160:HOH:O	1.77	0.85
1:E:98:PRO:HB3	1:F:98:PRO:HB3	1.61	0.82
1:F:96:ASP:O	1:F:124:GLN:HG2	1.81	0.81
1:D:1:MSE:CE	1:D:3:ALA:HB2	2.06	0.80
1:B:105:ALA:O	1:B:109:THR:HG23	1.83	0.78
1:D:20:GLU:O	1:D:24:GLN:NE2	2.16	0.78
1:E:20:GLU:O	1:E:24:GLN:HG2	1.83	0.78
1:D:52:ILE:HG13	1:D:83:MSE:HE3	1.67	0.76
1:F:18:ILE:CD1	1:F:100:VAL:HG22	2.17	0.75
1:C:10:HIS:NE2	1:C:72:SER:HB3	2.02	0.75
1:D:109:THR:OG1	1:D:122:GLU:OE2	2.06	0.74
1:A:36:LEU:HD22	1:A:50:LYS:HG2	1.68	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:49:MSE:HE1	1:C:52:ILE:HG13	1.71	0.72
1:D:1:MSE:HE1	1:D:65:LEU:HD11	1.70	0.72
1:C:96:ASP:O	1:C:124:GLN:HG2	1.91	0.71
1:F:18:ILE:HD11	1:F:100:VAL:HG22	1.71	0.70
1:B:50:LYS:NZ	2:B:149:HOH:O	2.21	0.70
1:E:28:SER:O	2:E:142:HOH:O	2.09	0.69
1:D:42:GLY:HA2	2:D:140:HOH:O	1.93	0.68
1:B:127:GLY:O	1:B:129:LYS:HE2	1.95	0.67
1:C:98:PRO:HB3	1:D:98:PRO:HB3	1.76	0.67
1:A:79:LEU:HD11	1:C:49:MSE:CE	2.19	0.65
1:A:79:LEU:HD13	1:C:49:MSE:HE3	1.72	0.65
1:D:48:PRO:HB2	1:D:83:MSE:HE1	1.78	0.65
1:F:87:ASP:OD1	2:F:162:HOH:O	2.14	0.64
1:C:96:ASP:O	1:C:124:GLN:CG	2.46	0.64
1:D:30:GLU:O	1:D:62:ARG:NH2	2.29	0.64
1:E:14:ILE:CD1	1:F:18:ILE:HD13	2.28	0.64
1:C:68:ALA:O	1:D:129:LYS:HE3	1.98	0.64
1:A:48:PRO:HG2	1:C:83:MSE:HE3	1.83	0.61
1:F:120:LEU:HB3	1:F:124:GLN:NE2	2.16	0.60
1:E:124:GLN:CG	2:E:160:HOH:O	2.43	0.60
1:F:120:LEU:HB3	1:F:124:GLN:HE22	1.64	0.60
1:C:41:ASP:OD2	1:E:32:THR:HG23	2.02	0.59
2:E:134:HOH:O	1:F:21:MSE:HE1	2.02	0.58
1:A:20:GLU:HG3	1:A:24:GLN:NE2	2.19	0.58
1:C:49:MSE:HE2	1:C:52:ILE:CG1	2.30	0.58
1:D:19:LYS:O	1:D:23:GLU:HG3	2.05	0.57
1:C:1:MSE:HB3	1:C:114:ASP:HA	1.86	0.57
1:E:29:GLU:HG3	1:E:31:ILE:HD13	1.86	0.56
1:A:79:LEU:HD12	1:C:49:MSE:HE3	1.80	0.56
1:F:88:GLN:HG3	2:F:162:HOH:O	2.06	0.55
1:C:1:MSE:HE3	1:C:119:ILE:HD12	1.87	0.55
1:E:14:ILE:HD11	1:F:18:ILE:HD13	1.89	0.55
1:A:106:SER:HA	1:A:109:THR:HG22	1.89	0.54
1:C:84:LEU:HB2	1:C:89:GLN:HG3	1.89	0.54
1:D:52:ILE:HG12	1:D:83:MSE:HE3	1.84	0.54
1:E:4:ASP:OD1	1:E:62:ARG:HD3	2.06	0.54
1:E:51:ILE:O	1:E:55:ILE:HG12	2.07	0.54
1:F:1:MSE:HE1	1:F:116:LEU:HB2	1.90	0.54
1:D:1:MSE:CE	1:D:3:ALA:CB	2.62	0.54
1:A:29:GLU:N	2:A:171:HOH:O	2.41	0.53
1:F:18:ILE:HD12	1:F:100:VAL:HG22	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:49:MSE:HE2	1:E:52:ILE:HD13	1.92	0.52
1:E:94:LEU:O	1:F:132:ASN:ND2	2.44	0.51
1:C:94:LEU:HB3	1:D:131:TRP:CE3	2.45	0.51
1:C:6:LEU:HD22	1:C:66:ILE:CD1	2.41	0.51
1:C:129:LYS:HE2	1:D:68:ALA:O	2.11	0.50
1:A:10:HIS:CD2	1:A:45:GLY:HA2	2.48	0.49
1:B:5:ILE:HD12	1:B:33:ILE:CD1	2.42	0.49
1:B:109:THR:HG21	1:B:122:GLU:HG3	1.94	0.49
1:C:74:VAL:O	1:C:78:GLU:HG3	2.13	0.48
1:A:94:LEU:HB3	1:B:131:TRP:CD2	2.48	0.48
1:D:106:SER:OG	1:D:123:ALA:HB2	2.14	0.48
1:D:1:MSE:HE2	1:D:110:ALA:HB1	1.94	0.48
1:E:93:HIS:CE1	1:E:116:LEU:HD21	2.49	0.47
1:E:39:THR:HG22	1:E:43:SER:H	1.80	0.47
1:F:9:SER:HB3	1:F:15:THR:HG23	1.96	0.47
1:E:96:ASP:O	1:E:124:GLN:HG3	2.14	0.47
1:A:109:THR:HG21	1:A:119:ILE:HA	1.96	0.46
1:D:52:ILE:HG12	1:D:83:MSE:CE	2.45	0.46
1:A:10:HIS:HE1	1:A:72:SER:CB	2.28	0.46
1:E:104:PHE:CE2	1:E:108:ILE:HD12	2.50	0.46
1:D:19:LYS:HE3	1:D:23:GLU:OE2	2.15	0.46
1:D:48:PRO:HB2	1:D:83:MSE:CE	2.46	0.45
1:A:131:TRP:CE3	1:B:94:LEU:HB3	2.52	0.45
1:C:49:MSE:HE2	1:C:49:MSE:HA	1.99	0.45
1:E:94:LEU:HB3	1:F:131:TRP:CE3	2.53	0.44
1:E:4:ASP:HB3	1:E:34:HIS:CD2	2.52	0.44
1:A:39:THR:HG22	1:A:43:SER:N	2.32	0.44
1:E:101:GLU:OE2	1:E:129:LYS:NZ	2.50	0.44
1:B:51:ILE:O	1:B:55:ILE:HG12	2.18	0.44
1:D:8:VAL:O	1:D:68:ALA:HA	2.18	0.44
1:E:8:VAL:O	1:E:68:ALA:HA	2.18	0.44
1:F:101:GLU:HG2	1:F:127:GLY:HA2	1.99	0.44
1:A:65:LEU:HD22	1:A:120:LEU:HD21	1.99	0.43
1:A:66:ILE:HG21	1:A:77:SER:HB3	2.01	0.43
1:C:68:ALA:O	1:D:129:LYS:CE	2.65	0.43
1:D:0:ALA:HB3	1:D:30:GLU:OE2	2.19	0.43
1:F:104:PHE:O	1:F:108:ILE:HG12	2.19	0.43
1:A:109:THR:CG2	1:A:119:ILE:HA	2.49	0.43
1:F:12:LYS:HE3	1:F:16:ASP:OD2	2.19	0.43
1:A:10:HIS:HE1	1:A:72:SER:HB3	1.82	0.42
1:A:98:PRO:HB3	1:B:98:PRO:HB3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:22:ILE:HG21	1:F:33:ILE:HG12	2.00	0.42
1:F:41:ASP:OD1	1:F:41:ASP:C	2.63	0.42
1:B:79:LEU:O	1:B:83:MSE:HG2	2.19	0.42
1:F:96:ASP:O	1:F:124:GLN:CG	2.59	0.42
1:F:106[B]:SER:OG	1:F:123:ALA:HB2	2.19	0.42
1:A:14:ILE:HD11	1:B:18:ILE:HA	2.01	0.42
1:B:101:GLU:HG2	1:B:129:LYS:NZ	2.35	0.41
1:C:25:MSE:HB3	2:C:173:HOH:O	2.21	0.41
1:E:39:THR:HG23	1:E:40:SER:N	2.35	0.41
1:F:8:VAL:O	1:F:68:ALA:HA	2.20	0.41
1:E:82:ASP:HB2	2:E:154:HOH:O	2.20	0.41
1:E:96:ASP:HB2	1:F:132:ASN:H	1.86	0.40
1:F:109:THR:OG1	1:F:122:GLU:HG2	2.21	0.40
1:D:46:SER:OG	1:D:73:ALA:HA	2.21	0.40
1:E:93:HIS:NE2	1:E:116:LEU:HD21	2.37	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	
1	A	123/135 (91%)	121 (98%)	2 (2%)	0	100	100
1	B	131/135 (97%)	125 (95%)	6 (5%)	0	100	100
1	C	132/135 (98%)	130 (98%)	2 (2%)	0	100	100
1	D	124/135 (92%)	122 (98%)	1 (1%)	1 (1%)	16	15
1	E	124/135 (92%)	121 (98%)	3 (2%)	0	100	100
1	F	134/135 (99%)	130 (97%)	4 (3%)	0	100	100
All	All	768/810 (95%)	749 (98%)	18 (2%)	1 (0%)	48	56

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	108	ILE

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	103/104 (99%)	99 (96%)	4 (4%)	28	37
1	B	107/104 (103%)	100 (94%)	7 (6%)	15	17
1	C	107/104 (103%)	106 (99%)	1 (1%)	70	82
1	D	101/104 (97%)	97 (96%)	4 (4%)	28	36
1	E	103/104 (99%)	91 (88%)	12 (12%)	5	4
1	F	109/104 (105%)	104 (95%)	5 (5%)	24	30
All	All	630/624 (101%)	597 (95%)	33 (5%)	20	25

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MSE
1	A	20	GLU
1	A	24	GLN
1	A	57	GLU
1	B	1	MSE
1	B	24	GLN
1	B	65	LEU
1	B	86	GLU
1	B	87	ASP
1	B	108	ILE
1	B	112	VAL
1	C	120	LEU
1	D	30	GLU
1	D	87	ASP
1	D	114	ASP
1	D	124	GLN
1	E	20	GLU
1	E	31	ILE

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Mol	Chain	Res	Type
1	E	39	THR
1	E	50	LYS
1	E	52	ILE
1	E	61	ASP
1	E	87	ASP
1	E	90	LYS
1	E	112	VAL
1	E	116	LEU
1	E	117	THR
1	E	122	GLU
1	F	10	HIS
1	F	46	SER
1	F	55	ILE
1	F	87	ASP
1	F	132	ASN

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	10	HIS
1	C	24	GLN
1	E	118	GLN
1	E	124	GLN
1	F	124	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 oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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 ⓘ

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	123/135 (91%)	0.29	1 (0%) 82 81	23, 29, 37, 58	0
1	B	126/135 (93%)	0.36	1 (0%) 82 81	14, 29, 36, 39	1 (0%)
1	C	128/135 (94%)	0.16	1 (0%) 82 81	21, 27, 36, 37	0
1	D	124/135 (91%)	0.56	9 (7%) 21 18	24, 30, 44, 55	0
1	E	124/135 (91%)	0.71	4 (3%) 50 47	21, 29, 42, 57	0
1	F	129/135 (95%)	0.20	4 (3%) 51 49	9, 28, 35, 44	1 (0%)
All	All	754/810 (93%)	0.38	20 (2%) 56 53	9, 29, 40, 58	2 (0%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	0	ALA	3.5
1	F	127	GLY	2.9
1	B	127	GLY	2.8
1	D	127	GLY	2.7
1	D	87	ASP	2.5
1	D	0	ALA	2.4
1	E	119	ILE	2.4
1	D	26	ASN	2.4
1	D	131	TRP	2.3
1	E	126	ALA	2.3
1	D	23	GLU	2.2
1	F	-2	SER	2.2
1	D	45	GLY	2.2
1	E	114	ASP	2.1
1	F	128	LYS	2.1
1	A	31	ILE	2.1
1	D	-1	ASN	2.1
1	F	41	ASP	2.1
1	C	127	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	116	LEU	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 oligosaccharides 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.