



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:38 AM UTC

PDB ID : 7XGE / pdb_00007xge
Title : Crystal structure of MCL-1 in complex with computationally designed inhibitor protein
Authors : Oh, B.-H.; Kim, S.
Deposited on : 2022-04-04
Resolution : 2.38 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

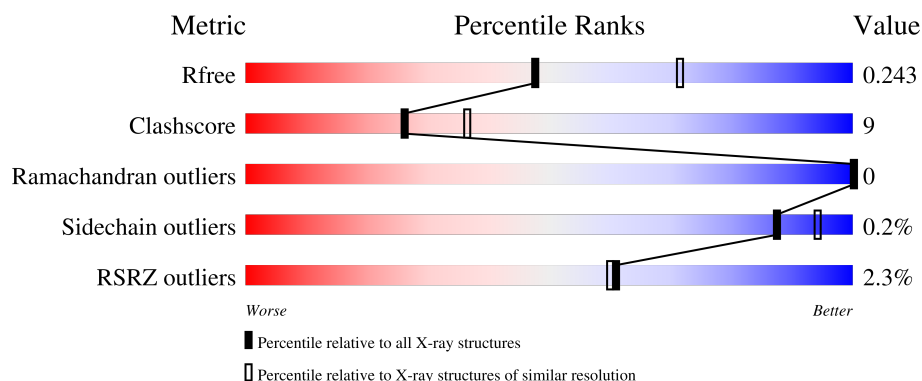
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

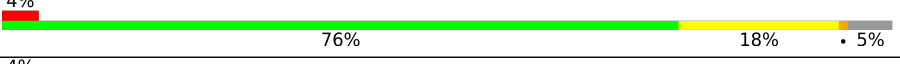
The reported resolution of this entry is 2.38 Å.

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	7164 (2.40-2.36)
Clashscore	190562	7722 (2.40-2.36)
Ramachandran outliers	187476	7626 (2.40-2.36)
Sidechain outliers	187428	7627 (2.40-2.36)
RSRZ outliers	180081	7170 (2.40-2.36)

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	169	
1	C	169	
1	E	169	
1	G	169	
2	B	156	

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Mol	Chain	Length	Quality of chain
2	D	156	% 77%15%8%
2	F	156	% 72%19%9%
2	H	156	% 69%21%11%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9956 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 BCL-xL and MCL-1 dual binder 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	168	Total	C	N	O	S	0	0	0
			1336	844	243	248	1			
1	C	161	Total	C	N	O	S	0	0	0
			1281	809	231	240	1			
1	E	151	Total	C	N	O	S	0	0	0
			1205	766	219	219	1			
1	G	164	Total	C	N	O	S	0	0	0
			1304	824	235	244	1			

- Molecule 2 is a protein called Induced myeloid leukemia cell differentiation protein Mcl-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	150	Total	C	N	O	S	0	0	0
			1208	759	223	222	4			
2	D	144	Total	C	N	O	S	0	0	0
			1162	731	213	215	3			
2	F	142	Total	C	N	O	S	0	0	0
			1145	721	210	211	3			
2	H	139	Total	C	N	O	S	0	0	0
			1128	711	207	207	3			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	3	MET	-	initiating methionine	UNP Q07820
B	154	SER	-	expression tag	UNP Q07820
B	155	ARG	-	expression tag	UNP Q07820
B	156	GLY	-	expression tag	UNP Q07820
B	157	LYS	-	expression tag	UNP Q07820
B	158	LEU	-	expression tag	UNP Q07820
D	3	MET	-	initiating methionine	UNP Q07820
D	154	SER	-	expression tag	UNP Q07820

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Chain	Residue	Modelled	Actual	Comment	Reference
D	155	ARG	-	expression tag	UNP Q07820
D	156	GLY	-	expression tag	UNP Q07820
D	157	LYS	-	expression tag	UNP Q07820
D	158	LEU	-	expression tag	UNP Q07820
F	3	MET	-	initiating methionine	UNP Q07820
F	154	SER	-	expression tag	UNP Q07820
F	155	ARG	-	expression tag	UNP Q07820
F	156	GLY	-	expression tag	UNP Q07820
F	157	LYS	-	expression tag	UNP Q07820
F	158	LEU	-	expression tag	UNP Q07820
H	3	MET	-	initiating methionine	UNP Q07820
H	154	SER	-	expression tag	UNP Q07820
H	155	ARG	-	expression tag	UNP Q07820
H	156	GLY	-	expression tag	UNP Q07820
H	157	LYS	-	expression tag	UNP Q07820
H	158	LEU	-	expression tag	UNP Q07820

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	25	Total O 25 25	0	0
3	B	23	Total O 23 23	0	0
3	C	34	Total O 34 34	0	0
3	D	21	Total O 21 21	0	0
3	E	22	Total O 22 22	0	0
3	F	19	Total O 19 19	0	0
3	G	24	Total O 24 24	0	0
3	H	19	Total O 19 19	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.

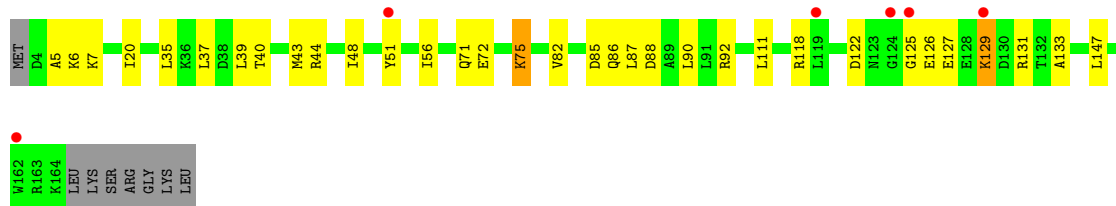
- Molecule 1: BCL-xL and MCL-1 dual binder 2

Chain A: 



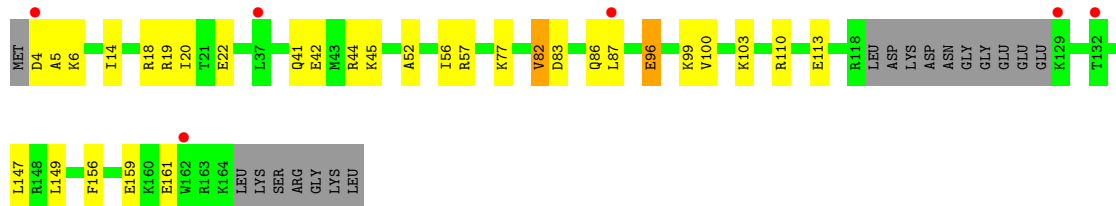
- Molecule 1: BCL-xL and MCL-1 dual binder 2

Chain C: 



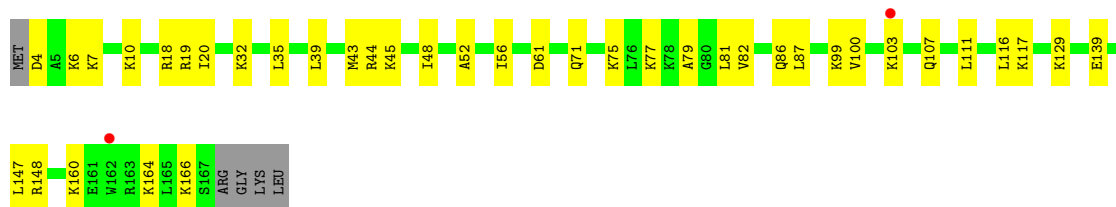
- Molecule 1: BCL-xL and MCL-1 dual binder 2

Chain E: 



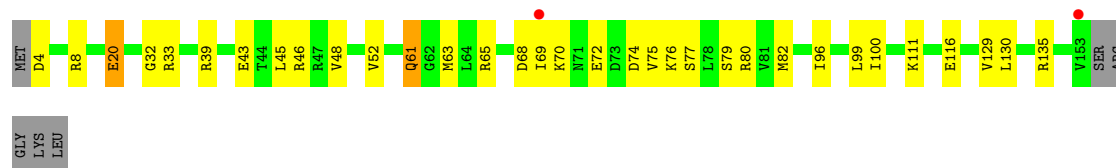
- Molecule 1: BCL-xL and MCL-1 dual binder 2

Chain G: 



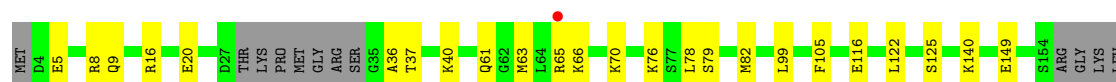
- Molecule 2: Induced myeloid leukemia cell differentiation protein Mcl-1

Chain B: 



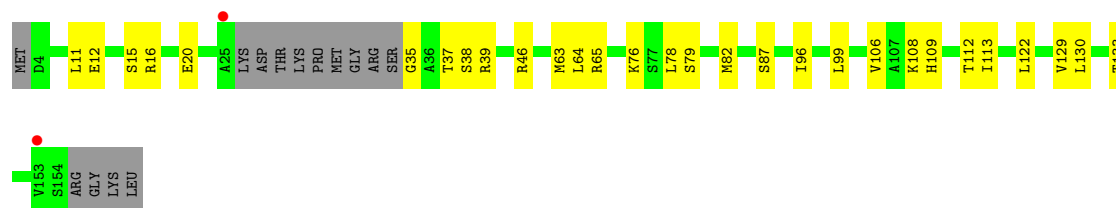
- Molecule 2: Induced myeloid leukemia cell differentiation protein Mcl-1

Chain D: 



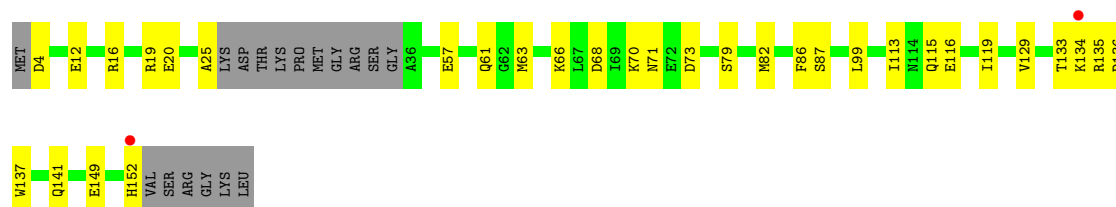
- Molecule 2: Induced myeloid leukemia cell differentiation protein Mcl-1

Chain F: 



- Molecule 2: Induced myeloid leukemia cell differentiation protein Mcl-1

Chain H: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	67.99Å 74.77Å 92.16Å 70.83° 89.59° 67.58°	Depositor
Resolution (Å)	29.46 – 2.38 29.46 – 2.38	Depositor EDS
% Data completeness (in resolution range)	97.3 (29.46-2.38) 97.3 (29.46-2.38)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.93 (at 2.39Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.202 , 0.240 0.203 , 0.243	Depositor DCC
R_{free} test set	3065 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å ²)	53.5	Xtriage
Anisotropy	0.264	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 49.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9956	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 7.81% 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	0.24	0/1345	0.52	1/1795 (0.1%)
1	C	0.31	0/1290	0.62	4/1724 (0.2%)
1	E	0.35	0/1213	0.66	2/1620 (0.1%)
1	G	0.25	0/1313	0.52	0/1754
2	B	0.28	0/1228	0.57	3/1651 (0.2%)
2	D	0.33	0/1180	0.76	7/1586 (0.4%)
2	F	0.22	0/1163	0.52	1/1564 (0.1%)
2	H	0.24	0/1146	0.47	0/1541
All	All	0.28	0/9878	0.59	18/13235 (0.1%)

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	75	LYS	CB-CG-CD	9.28	132.65	111.30
2	D	65	ARG	CB-CG-CD	-8.36	92.07	111.30
1	A	75	LYS	CA-CB-CG	8.13	130.37	114.10
2	D	65	ARG	NE-CZ-NH2	6.86	125.37	119.20
1	E	96	GLU	CA-CB-CG	6.63	127.36	114.10
2	D	66	LYS	CA-CB-CG	6.43	126.96	114.10
2	D	65	ARG	CG-CD-NE	6.00	125.21	112.00
2	D	65	ARG	NE-CZ-NH1	-5.97	115.53	121.50
2	D	65	ARG	CA-CB-CG	5.90	125.90	114.10
2	B	20	GLU	CB-CG-CD	-5.84	102.68	112.60
2	B	20	GLU	CA-CB-CG	5.79	125.68	114.10
2	F	37	THR	OG1-CB-CG2	5.27	119.84	109.30
1	C	129	LYS	CB-CG-CD	5.26	123.40	111.30
1	E	161	GLU	N-CA-CB	5.23	118.32	110.53
2	B	69	ILE	CA-CB-CG1	5.06	119.00	110.40
1	C	87	LEU	CA-C-N	5.03	129.11	120.72
1	C	87	LEU	C-N-CA	5.03	129.11	120.72
2	D	66	LYS	CB-CG-CD	5.01	122.82	111.30

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	1336	0	1417	18	0
1	C	1281	0	1348	36	1
1	E	1205	0	1285	28	0
1	G	1304	0	1376	30	1
2	B	1208	0	1220	22	0
2	D	1162	0	1167	19	0
2	F	1145	0	1150	22	0
2	H	1128	0	1133	26	0
3	A	25	0	0	0	0
3	B	23	0	0	0	0
3	C	34	0	0	1	0
3	D	21	0	0	0	0
3	E	22	0	0	0	0
3	F	19	0	0	0	0
3	G	24	0	0	1	0
3	H	19	0	0	0	0
All	All	9956	0	10096	187	1

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 9.

All (187) 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)
2:F:76:LYS:NZ	2:F:79:SER:OG	1.76	1.18
1:C:82:VAL:HG23	1:C:86:GLN:HB2	1.20	1.12
1:C:126:GLU:HA	1:C:129:LYS:HD2	1.45	0.97
2:F:20:GLU:OE1	2:F:46:ARG:NH1	1.99	0.94
1:C:82:VAL:CG2	1:C:86:GLN:HB2	2.00	0.91
1:C:71:GLN:HG2	1:C:75:LYS:HE3	1.51	0.90
2:B:111:LYS:NZ	2:B:116:GLU:OE1	2.07	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:82:VAL:HG23	1:G:86:GLN:HB2	1.55	0.87
1:C:125:GLY:O	1:C:129:LYS:HG2	1.78	0.83
1:C:37:LEU:HD21	1:C:129:LYS:NZ	1.95	0.81
1:A:6:LYS:HD2	1:A:159:GLU:HG2	1.63	0.80
1:E:6:LYS:HE2	1:E:159:GLU:HG2	1.63	0.80
1:C:35:LEU:HD22	1:C:129:LYS:HB2	1.63	0.79
2:B:76:LYS:HE3	2:B:80:ARG:HH12	1.51	0.76
1:C:37:LEU:HD21	1:C:129:LYS:HZ2	1.48	0.76
1:E:113:GLU:OE1	2:F:65:ARG:NH1	2.19	0.75
1:G:82:VAL:HG23	1:G:86:GLN:CB	2.16	0.74
2:F:12:GLU:O	2:F:16:ARG:HG3	1.87	0.73
1:E:83:ASP:H	1:E:86:GLN:HE21	1.35	0.73
2:B:20:GLU:OE1	2:B:46:ARG:NE	2.23	0.69
1:C:85:ASP:HA	1:C:88:ASP:OD1	1.92	0.69
1:G:35:LEU:HD22	1:G:129:LYS:HG2	1.75	0.69
1:G:52:ALA:HB1	2:H:63:MET:HE3	1.76	0.68
1:E:18:ARG:HD2	1:E:19:ARG:HD3	1.77	0.67
2:F:35:GLY:N	2:F:38:SER:HG	1.93	0.66
1:E:99:LYS:NZ	1:E:100:VAL:HG13	2.10	0.66
2:D:36:ALA:O	2:D:40:LYS:HG3	1.97	0.65
1:C:20:ILE:HG13	1:C:147:LEU:HD22	1.77	0.64
2:H:57:GLU:O	2:H:61:GLN:HB2	1.98	0.64
1:A:39:LEU:HD23	1:A:43:MET:HE2	1.80	0.64
1:C:85:ASP:O	1:C:88:ASP:OD1	2.16	0.64
2:D:61:GLN:HG2	2:D:105:PHE:CZ	2.33	0.64
1:C:85:ASP:C	1:C:88:ASP:OD1	2.42	0.63
1:A:68:GLN:O	1:A:71:GLN:HG2	1.98	0.62
2:F:76:LYS:NZ	2:F:79:SER:HG	1.97	0.61
2:D:82:MET:HG2	2:D:99:LEU:HD22	1.84	0.60
2:D:76:LYS:O	2:D:79:SER:HB2	2.02	0.60
1:G:18:ARG:NH1	1:G:148:ARG:HH22	2.01	0.59
2:H:149:GLU:O	2:H:152:HIS:ND1	2.36	0.58
1:G:20:ILE:HG13	1:G:147:LEU:HD22	1.86	0.58
1:E:82:VAL:HG22	1:E:86:GLN:HB2	1.85	0.58
1:C:72:GLU:O	1:C:75:LYS:HB2	2.04	0.57
2:B:82:MET:HG2	2:B:99:LEU:HD22	1.87	0.57
1:C:71:GLN:HG2	1:C:75:LYS:CE	2.29	0.57
1:G:39:LEU:HD23	1:G:43:MET:HE2	1.87	0.57
1:C:56:ILE:HG13	2:D:63:MET:HE1	1.87	0.57
2:H:68:ASP:CG	2:H:70:LYS:HE3	2.30	0.56
1:C:147:LEU:O	1:C:147:LEU:HG	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:149:GLU:O	2:D:149:GLU:HG3	2.03	0.56
1:E:42:GLU:OE1	1:E:42:GLU:N	2.34	0.56
2:B:61:GLN:O	2:B:65:ARG:HG3	2.05	0.56
1:A:160:LYS:HD3	1:A:163:ARG:HH21	1.71	0.56
2:H:16:ARG:HG2	2:H:19:ARG:NH2	2.22	0.55
1:C:72:GLU:HA	1:C:75:LYS:CG	2.37	0.55
1:G:6:LYS:HG3	1:G:7:LYS:N	2.21	0.55
1:E:41:GLN:O	1:E:45:LYS:HG2	2.06	0.55
2:F:82:MET:HG2	2:F:99:LEU:HD22	1.88	0.55
2:H:86:PHE:CG	2:H:134:LYS:HE2	2.42	0.54
2:D:61:GLN:HG2	2:D:105:PHE:HZ	1.71	0.54
1:E:99:LYS:O	1:E:103:LYS:HG3	2.09	0.53
1:G:111:LEU:HD21	1:G:139:GLU:HB3	1.90	0.53
1:G:56:ILE:HG13	2:H:63:MET:HE1	1.91	0.52
2:F:79:SER:HB2	2:F:129:VAL:HG21	1.91	0.52
2:H:57:GLU:HG2	2:H:61:GLN:HE21	1.75	0.52
1:C:85:ASP:CA	1:C:88:ASP:OD1	2.56	0.52
1:E:99:LYS:HD2	1:E:100:VAL:N	2.25	0.51
2:H:82:MET:HG2	2:H:99:LEU:HD22	1.93	0.51
2:D:9:GLN:NE2	2:D:37:THR:OG1	2.43	0.51
2:B:45:LEU:HD21	2:B:100:ILE:HG21	1.93	0.51
1:C:126:GLU:HA	1:C:129:LYS:CD	2.30	0.51
1:E:96:GLU:CD	1:E:99:LYS:HE3	2.35	0.51
1:E:99:LYS:HZ2	1:E:100:VAL:HG13	1.75	0.51
1:C:37:LEU:CD2	1:C:129:LYS:NZ	2.71	0.50
1:A:160:LYS:HD3	1:A:163:ARG:NH2	2.27	0.50
1:G:44:ARG:O	1:G:48:ILE:HG12	2.10	0.50
1:G:18:ARG:HH12	1:G:148:ARG:HH22	1.59	0.50
2:D:5:GLU:OE2	2:D:8:ARG:NH1	2.39	0.50
2:H:113:ILE:HD12	2:H:115:GLN:HG3	1.93	0.50
1:A:20:ILE:HG13	1:A:147:LEU:HD22	1.93	0.49
1:A:68:GLN:HG2	1:A:71:GLN:NE2	2.27	0.49
2:B:20:GLU:OE1	2:B:46:ARG:NH2	2.44	0.49
2:B:96:ILE:HG12	2:B:130:LEU:HD21	1.94	0.49
2:H:12:GLU:O	2:H:16:ARG:HG3	2.10	0.49
2:D:78:LEU:HD23	2:D:122:LEU:HD11	1.94	0.49
2:F:78:LEU:HD23	2:F:122:LEU:HD11	1.94	0.49
2:F:96:ILE:HG12	2:F:130:LEU:HD11	1.95	0.49
2:D:140:LYS:O	2:D:140:LYS:HD3	2.11	0.49
1:A:49:GLU:OE2	2:B:77:SER:OG	2.25	0.49
1:E:147:LEU:O	1:E:147:LEU:HG	2.10	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:4:ASP:OD2	2:B:135:ARG:NH2	2.39	0.49
1:C:35:LEU:HD11	1:C:133:ALA:HB2	1.95	0.49
1:C:118:ARG:O	1:C:122:ASP:HB2	2.12	0.49
1:G:82:VAL:CG2	1:G:86:GLN:HB2	2.34	0.49
2:B:39:ARG:O	2:B:43:GLU:HG3	2.14	0.48
1:A:122:ASP:OD1	1:A:128:GLU:HG3	2.13	0.48
2:F:64:LEU:HD13	2:F:106:VAL:HG22	1.96	0.48
2:H:79:SER:HB2	2:H:129:VAL:HG21	1.96	0.48
1:C:72:GLU:HA	1:C:75:LYS:HG2	1.95	0.48
1:G:45:LYS:HE2	1:G:45:LYS:HB2	1.54	0.48
1:G:99:LYS:O	1:G:103:LYS:HG2	2.14	0.48
2:D:5:GLU:CD	2:D:8:ARG:NH1	2.72	0.48
1:G:77:LYS:HB2	1:G:87:LEU:HD13	1.95	0.48
2:D:79:SER:OG	2:D:125:SER:OG	2.32	0.47
1:E:96:GLU:O	1:E:100:VAL:HG22	2.14	0.47
2:B:111:LYS:CE	2:B:116:GLU:OE1	2.63	0.47
1:G:116:LEU:HD13	2:H:66:LYS:HD2	1.95	0.47
1:A:72:GLU:O	1:A:75:LYS:HB3	2.14	0.47
1:C:39:LEU:HD12	1:C:43:MET:HE2	1.97	0.47
1:E:110:ARG:NH1	1:E:113:GLU:HB3	2.30	0.47
2:F:35:GLY:O	2:F:39:ARG:HG3	2.14	0.47
2:B:72:GLU:O	2:B:75:VAL:HG12	2.15	0.46
1:G:166:LYS:HE3	1:G:166:LYS:HA	1.96	0.46
2:B:70:LYS:HD2	2:B:74:ASP:OD2	2.14	0.46
2:H:4:ASP:OD2	2:H:135:ARG:NH2	2.44	0.46
1:E:56:ILE:HG13	2:F:63:MET:HE1	1.96	0.46
1:G:6:LYS:HD3	1:G:10:LYS:HZ2	1.79	0.46
1:C:40:THR:HG23	1:C:43:MET:H	1.81	0.46
1:E:20:ILE:HG13	1:E:147:LEU:HD22	1.96	0.46
1:E:4:ASP:OD1	1:E:5:ALA:N	2.49	0.46
2:H:71:ASN:OD1	2:H:73:ASP:HB2	2.16	0.45
1:C:88:ASP:O	1:C:92:ARG:HG3	2.16	0.45
2:B:48:VAL:O	2:B:52:VAL:HG23	2.17	0.45
1:E:149:LEU:HD12	1:E:149:LEU:HA	1.79	0.45
2:F:78:LEU:O	2:F:82:MET:HG3	2.16	0.45
1:G:19:ARG:HD2	1:G:61:ASP:HB3	1.99	0.45
1:C:126:GLU:O	1:C:129:LYS:HG3	2.17	0.45
2:H:137:TRP:CH2	2:H:141:GLN:HG3	2.52	0.45
2:D:16:ARG:O	2:D:20:GLU:HB3	2.18	0.44
2:D:5:GLU:OE1	2:D:8:ARG:NH1	2.51	0.44
1:A:7:LYS:HA	1:A:7:LYS:HD3	1.72	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:116:GLU:HA	2:H:119:ILE:HG13	1.98	0.44
1:A:134:LYS:HG3	1:E:156:PHE:HZ	1.82	0.44
2:D:5:GLU:CD	2:D:8:ARG:HH12	2.26	0.44
1:A:120:ASP:OD2	1:A:121:LYS:HE3	2.18	0.44
2:D:116:GLU:H	2:D:116:GLU:CD	2.26	0.44
1:G:4:ASP:HB2	1:G:6:LYS:HG2	2.00	0.44
1:G:32:LYS:HB3	1:G:32:LYS:HE2	1.84	0.44
1:A:68:GLN:O	1:A:72:GLU:HG2	2.17	0.43
1:C:129:LYS:HG2	1:C:129:LYS:H	1.61	0.43
2:B:8:ARG:NH1	2:B:32:GLY:HA3	2.33	0.43
1:E:52:ALA:HB1	2:F:63:MET:HE3	2.00	0.43
1:A:44:ARG:O	1:A:48:ILE:HG12	2.17	0.43
1:C:37:LEU:CD2	1:C:129:LYS:HZ3	2.31	0.43
1:G:99:LYS:HG3	1:G:100:VAL:N	2.32	0.43
1:G:79:ALA:HB3	1:G:81:LEU:HD12	2.00	0.43
2:B:33:ARG:NE	2:B:33:ARG:HA	2.33	0.43
2:B:68:ASP:CG	2:B:70:LYS:HE3	2.44	0.43
1:C:5:ALA:HB1	1:C:90:LEU:HD22	2.00	0.43
1:E:99:LYS:HZ3	1:E:100:VAL:HG13	1.82	0.43
2:F:11:LEU:O	2:F:15:SER:HB2	2.19	0.43
2:F:108:LYS:O	2:F:112:THR:HG23	2.19	0.43
1:A:34:ALA:HB1	1:A:119:LEU:HD21	2.00	0.42
2:B:76:LYS:H	2:B:76:LYS:HG2	1.65	0.42
1:C:6:LYS:HG3	1:C:7:LYS:N	2.34	0.42
2:F:87:SER:HA	2:H:133:THR:O	2.19	0.42
2:F:16:ARG:O	2:F:20:GLU:HG2	2.20	0.42
2:D:61:GLN:CG	2:D:105:PHE:HZ	2.32	0.42
1:C:111:LEU:HD23	1:C:111:LEU:HA	1.93	0.42
1:G:71:GLN:O	1:G:75:LYS:HG3	2.19	0.42
1:C:72:GLU:HA	1:C:75:LYS:HG3	2.02	0.42
1:C:127:GLU:CD	1:C:131:ARG:NH2	2.78	0.42
1:E:57:ARG:NH2	2:H:136:ASP:OD2	2.37	0.42
2:H:68:ASP:OD2	2:H:70:LYS:HE3	2.20	0.42
1:C:44:ARG:O	1:C:48:ILE:HG12	2.20	0.41
2:B:79:SER:HB2	2:B:129:VAL:HG21	2.02	0.41
1:E:44:ARG:HH11	1:E:45:LYS:NZ	2.19	0.41
2:H:86:PHE:CD2	2:H:134:LYS:HE2	2.55	0.41
2:F:109:HIS:O	2:F:113:ILE:HG12	2.20	0.41
1:C:51:TYR:OH	3:C:201:HOH:O	2.17	0.41
1:E:77:LYS:HB2	1:E:87:LEU:HD13	2.03	0.41
2:F:133:THR:O	2:H:87:SER:HA	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:160:LYS:HG2	1:G:164:LYS:HE2	2.03	0.41
2:H:20:GLU:HG3	2:H:25:ALA:C	2.46	0.41
1:E:44:ARG:HD3	1:E:45:LYS:HZ1	1.85	0.41
2:F:76:LYS:HE3	2:F:76:LYS:HB3	1.76	0.41
1:G:18:ARG:NH1	3:G:201:HOH:O	2.24	0.41
2:D:70:LYS:HE3	2:D:70:LYS:HB2	1.75	0.41
1:E:14:ILE:HD13	1:E:14:ILE:HA	1.89	0.41
2:H:133:THR:O	2:H:134:LYS:HG3	2.21	0.41
1:A:52:ALA:HB1	2:B:63:MET:HE3	2.03	0.41
2:B:20:GLU:OE1	2:B:46:ARG:CZ	2.69	0.40
1:A:32:LYS:NZ	1:A:137:GLU:OE2	2.36	0.40
1:E:18:ARG:HD3	1:E:22:GLU:OE2	2.21	0.40
1:G:103:LYS:O	1:G:107:GLN:HG3	2.21	0.40
1:G:117:LYS:O	1:G:117:LYS:HD2	2.21	0.40
2:H:16:ARG:HG2	2:H:19:ARG:HH22	1.86	0.40
1:G:48:ILE:HD12	2:H:66:LYS:HB3	2.04	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:131:ARG:NH2	1:G:7:LYS:NZ[1_556]	2.13	0.07

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	166/169 (98%)	164 (99%)	2 (1%)	0	100	100
1	C	159/169 (94%)	157 (99%)	2 (1%)	0	100	100
1	E	147/169 (87%)	146 (99%)	1 (1%)	0	100	100
1	G	162/169 (96%)	160 (99%)	2 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	148/156 (95%)	142 (96%)	6 (4%)	0	100	100
2	D	140/156 (90%)	138 (99%)	2 (1%)	0	100	100
2	F	138/156 (88%)	136 (99%)	2 (1%)	0	100	100
2	H	135/156 (86%)	132 (98%)	3 (2%)	0	100	100
All	All	1195/1300 (92%)	1175 (98%)	20 (2%)	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	133/134 (99%)	133 (100%)	0	100	100
1	C	127/134 (95%)	127 (100%)	0	100	100
1	E	119/134 (89%)	118 (99%)	1 (1%)	73	85
1	G	130/134 (97%)	130 (100%)	0	100	100
2	B	131/136 (96%)	130 (99%)	1 (1%)	73	85
2	D	126/136 (93%)	126 (100%)	0	100	100
2	F	124/136 (91%)	124 (100%)	0	100	100
2	H	122/136 (90%)	122 (100%)	0	100	100
All	All	1012/1080 (94%)	1010 (100%)	2 (0%)	87	94

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

Mol	Chain	Res	Type
2	B	61	GLN
1	E	82	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	GLN
1	A	71	GLN
1	A	98	GLN
2	D	9	GLN
2	D	71	ASN
1	E	86	GLN
2	F	71	ASN
2	F	109	HIS
1	G	107	GLN
2	H	61	GLN

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 oligosaccharides 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(Å ²)	Q < 0.9
1	A	168/169 (99%)	0.23	7 (4%)	40	40	42, 63, 107, 156	0
1	C	161/169 (95%)	0.21	6 (3%)	45	45	42, 67, 130, 157	0
1	E	151/169 (89%)	0.28	6 (3%)	42	42	43, 66, 121, 139	0
1	G	164/169 (97%)	0.12	2 (1%)	76	76	41, 64, 107, 126	0
2	B	150/156 (96%)	0.09	2 (1%)	75	74	40, 64, 114, 148	0
2	D	144/156 (92%)	0.11	1 (0%)	84	83	42, 63, 93, 137	0
2	F	142/156 (91%)	0.13	2 (1%)	73	72	44, 67, 101, 120	0
2	H	139/156 (89%)	0.10	2 (1%)	73	72	42, 69, 102, 116	0
All	All	1219/1300 (93%)	0.16	28 (2%)	61	60	40, 65, 113, 157	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	162	TRP	4.9
2	B	153	VAL	4.7
1	C	162	TRP	4.3
1	A	119	LEU	3.8
1	G	162	TRP	3.3
1	C	119	LEU	2.9
1	A	124	GLY	2.8
2	F	25	ALA	2.7
2	F	153	VAL	2.7
1	C	51	TYR	2.6
1	E	87	LEU	2.6
2	H	152	HIS	2.6
1	E	129	LYS	2.6
1	C	125	GLY	2.5
1	A	171	LEU	2.4
2	D	65	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	128	GLU	2.3
1	A	120	ASP	2.3
1	A	35	LEU	2.3
2	B	69	ILE	2.2
1	G	103	LYS	2.2
1	E	132	THR	2.2
1	C	129	LYS	2.1
1	A	162	TRP	2.1
1	C	124	GLY	2.1
2	H	134	LYS	2.1
1	E	4	ASP	2.0
1	E	37	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.