



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:38 AM GMT

PDB ID : 3BIM
Title : Crystal structure of the BCL6 BTB domain dimer in complex with the BCOR
BBD corepressor peptide
Authors : Prive, G.G.; Ghetu, A.F.
Deposited on : 2007-11-30
Resolution : 2.60 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
<http://wwpdb.org/validation/2016/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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

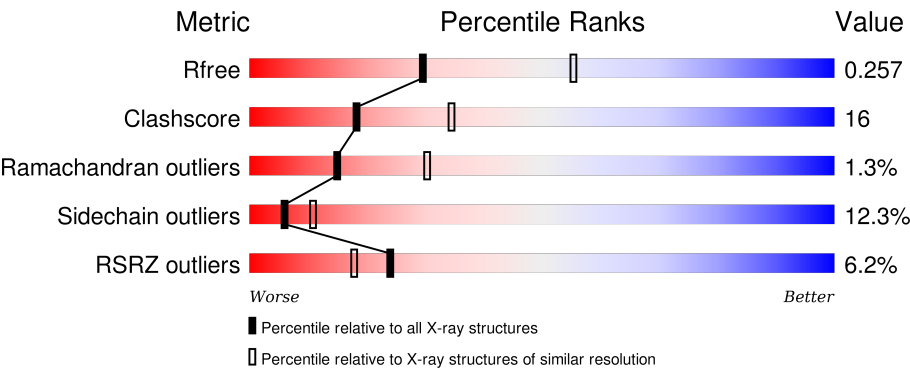
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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}	91344	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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. 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	127	
1	B	127	
1	C	127	
1	D	127	
1	E	127	

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Mol	Chain	Length	Quality of chain
1	F	127	
1	G	127	
1	H	127	
2	I	19	
2	J	19	
2	K	19	
2	L	19	
2	M	19	
2	N	19	
2	O	19	
2	P	19	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9091 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 B-cell lymphoma 6 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	125	Total	C	N	O	S	0	0	0
			1008	636	178	186	8			
1	B	124	Total	C	N	O	S	0	0	0
			998	631	177	182	8			
1	C	125	Total	C	N	O	S	0	0	0
			1008	636	178	186	8			
1	D	124	Total	C	N	O	S	0	0	0
			998	631	177	182	8			
1	E	123	Total	C	N	O	S	0	0	0
			993	628	176	181	8			
1	F	122	Total	C	N	O	S	0	0	0
			987	625	175	179	8			
1	G	122	Total	C	N	O	S	0	0	0
			985	624	175	178	8			
1	H	122	Total	C	N	O	S	0	0	0
			987	625	175	179	8			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	3	GLY	-	EXPRESSION TAG	UNP P41182
A	4	SER	-	EXPRESSION TAG	UNP P41182
A	8	GLN	CYS	ENGINEERED	UNP P41182
A	67	ARG	CYS	ENGINEERED	UNP P41182
A	84	ASN	CYS	ENGINEERED	UNP P41182
B	3	GLY	-	EXPRESSION TAG	UNP P41182
B	4	SER	-	EXPRESSION TAG	UNP P41182
B	8	GLN	CYS	ENGINEERED	UNP P41182
B	67	ARG	CYS	ENGINEERED	UNP P41182
B	84	ASN	CYS	ENGINEERED	UNP P41182
C	3	GLY	-	EXPRESSION TAG	UNP P41182
C	4	SER	-	EXPRESSION TAG	UNP P41182
C	8	GLN	CYS	ENGINEERED	UNP P41182

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Chain	Residue	Modelled	Actual	Comment	Reference
C	67	ARG	CYS	ENGINEERED	UNP P41182
C	84	ASN	CYS	ENGINEERED	UNP P41182
D	3	GLY	-	EXPRESSION TAG	UNP P41182
D	4	SER	-	EXPRESSION TAG	UNP P41182
D	8	GLN	CYS	ENGINEERED	UNP P41182
D	67	ARG	CYS	ENGINEERED	UNP P41182
D	84	ASN	CYS	ENGINEERED	UNP P41182
E	3	GLY	-	EXPRESSION TAG	UNP P41182
E	4	SER	-	EXPRESSION TAG	UNP P41182
E	8	GLN	CYS	ENGINEERED	UNP P41182
E	67	ARG	CYS	ENGINEERED	UNP P41182
E	84	ASN	CYS	ENGINEERED	UNP P41182
F	3	GLY	-	EXPRESSION TAG	UNP P41182
F	4	SER	-	EXPRESSION TAG	UNP P41182
F	8	GLN	CYS	ENGINEERED	UNP P41182
F	67	ARG	CYS	ENGINEERED	UNP P41182
F	84	ASN	CYS	ENGINEERED	UNP P41182
G	3	GLY	-	EXPRESSION TAG	UNP P41182
G	4	SER	-	EXPRESSION TAG	UNP P41182
G	8	GLN	CYS	ENGINEERED	UNP P41182
G	67	ARG	CYS	ENGINEERED	UNP P41182
G	84	ASN	CYS	ENGINEERED	UNP P41182
H	3	GLY	-	EXPRESSION TAG	UNP P41182
H	4	SER	-	EXPRESSION TAG	UNP P41182
H	8	GLN	CYS	ENGINEERED	UNP P41182
H	67	ARG	CYS	ENGINEERED	UNP P41182
H	84	ASN	CYS	ENGINEERED	UNP P41182

- Molecule 2 is a protein called BCL-6 corepressor.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	I	17	Total	C	N	O	0	0	0
			126	80	21	25			
2	J	17	Total	C	N	O	0	0	0
			126	80	21	25			
2	K	17	Total	C	N	O	0	0	0
			126	80	21	25			
2	L	16	Total	C	N	O	0	0	0
			115	74	17	24			
2	M	17	Total	C	N	O	0	0	0
			126	80	21	25			
2	N	16	Total	C	N	O	0	0	0
			115	74	17	24			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	O	17	Total	C	N	O	0	0	0
			126	80	21	25			
2	P	17	Total	C	N	O	0	0	0
			126	80	21	25			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
I	496	GLY	-	EXPRESSION TAG	UNP Q6W2J9
I	497	SER	-	EXPRESSION TAG	UNP Q6W2J9
J	496	GLY	-	EXPRESSION TAG	UNP Q6W2J9
J	497	SER	-	EXPRESSION TAG	UNP Q6W2J9
K	496	GLY	-	EXPRESSION TAG	UNP Q6W2J9
K	497	SER	-	EXPRESSION TAG	UNP Q6W2J9
L	496	GLY	-	EXPRESSION TAG	UNP Q6W2J9
L	497	SER	-	EXPRESSION TAG	UNP Q6W2J9
M	496	GLY	-	EXPRESSION TAG	UNP Q6W2J9
M	497	SER	-	EXPRESSION TAG	UNP Q6W2J9
N	496	GLY	-	EXPRESSION TAG	UNP Q6W2J9
N	497	SER	-	EXPRESSION TAG	UNP Q6W2J9
O	496	GLY	-	EXPRESSION TAG	UNP Q6W2J9
O	497	SER	-	EXPRESSION TAG	UNP Q6W2J9
P	496	GLY	-	EXPRESSION TAG	UNP Q6W2J9
P	497	SER	-	EXPRESSION TAG	UNP Q6W2J9

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	26	Total	O	0	0
			26	26		
3	B	23	Total	O	0	0
			23	23		
3	C	11	Total	O	0	0
			11	11		
3	D	23	Total	O	0	0
			23	23		
3	E	12	Total	O	0	0
			12	12		
3	F	15	Total	O	0	0
			15	15		
3	G	15	Total	O	0	0
			15	15		

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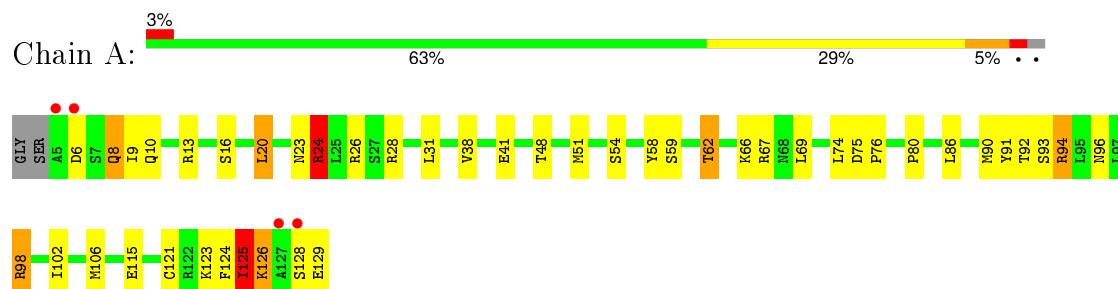
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	H	7	Total 7	O 7	0	0
3	I	1	Total 1	O 1	0	0
3	J	3	Total 3	O 3	0	0
3	K	1	Total 1	O 1	0	0
3	L	1	Total 1	O 1	0	0
3	M	1	Total 1	O 1	0	0
3	N	1	Total 1	O 1	0	0
3	O	1	Total 1	O 1	0	0

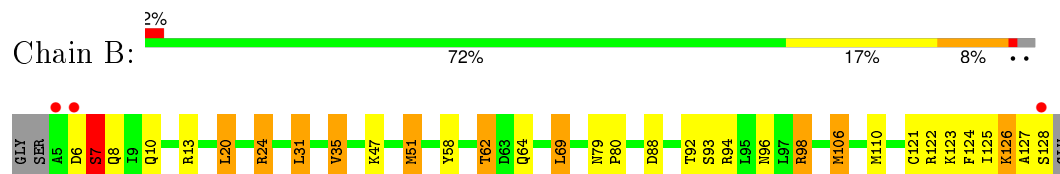
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

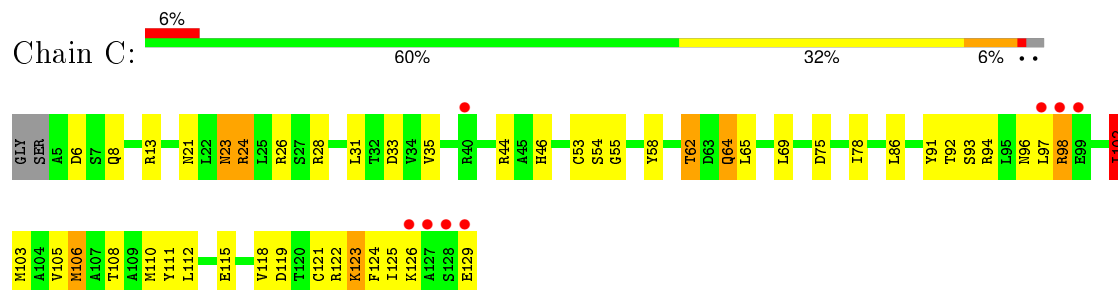
- Molecule 1: B-cell lymphoma 6 protein



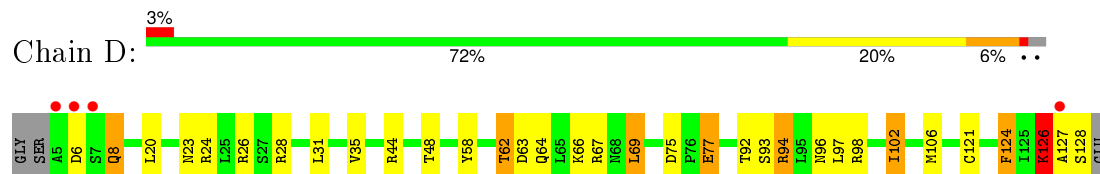
- Molecule 1: B-cell lymphoma 6 protein



- Molecule 1: B-cell lymphoma 6 protein

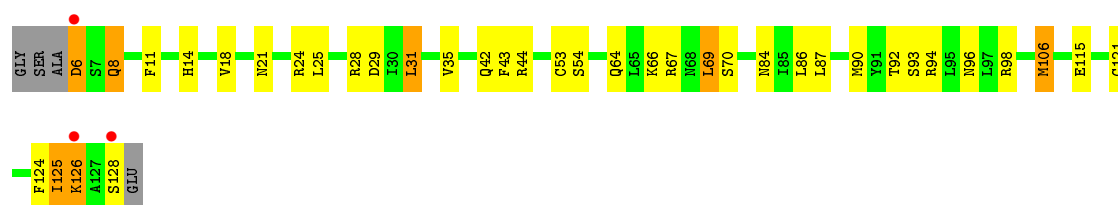


- Molecule 1: B-cell lymphoma 6 protein

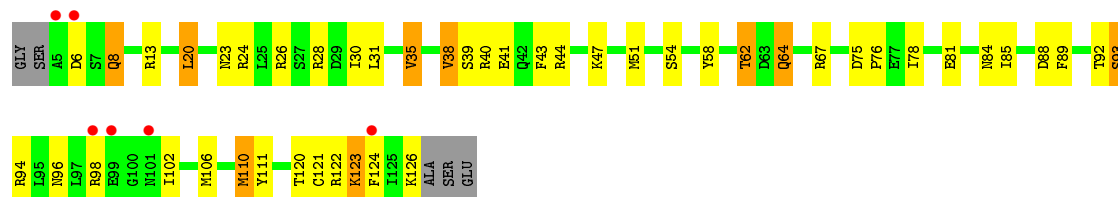


- Molecule 1: B-cell lymphoma 6 protein

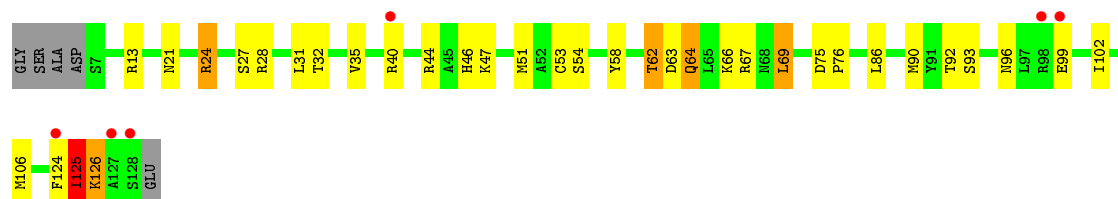




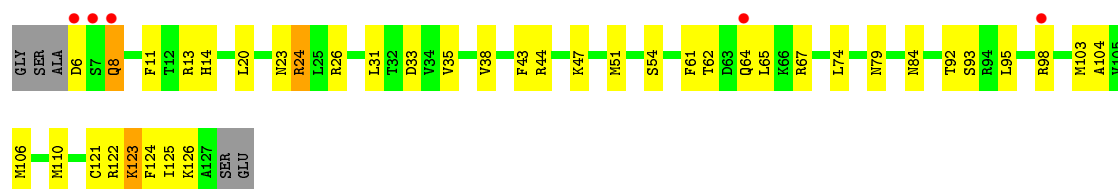
- Molecule 1: B-cell lymphoma 6 protein



- Molecule 1: B-cell lymphoma 6 protein



- Molecule 1: B-cell lymphoma 6 protein

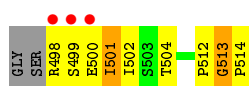


- Molecule 2: BCL-6 corepressor

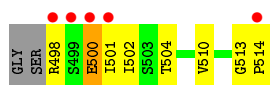


- Molecule 2: BCL-6 corepressor

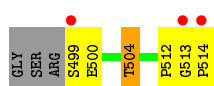




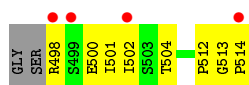
- Molecule 2: BCL-6 corepressor



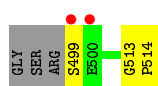
- Molecule 2: BCL-6 corepressor



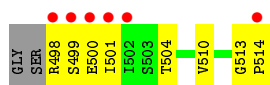
- Molecule 2: BCL-6 corepressor



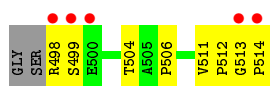
- Molecule 2: BCL-6 corepressor



- Molecule 2: BCL-6 corepressor



- Molecule 2: BCL-6 corepressor



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	150.55Å 150.55Å 312.25Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.70 – 2.60 48.68 – 2.40	Depositor EDS
% Data completeness (in resolution range)	97.1 (48.70-2.60) 98.3 (48.68-2.40)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.219 , 0.260 0.216 , 0.257	Depositor DCC
R_{free} test set	3159 reflections (5.12%)	DCC
Wilson B-factor (Å ²)	62.7	Xtriage
Anisotropy	0.315	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 55.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 80574 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9091	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.09% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.88	1/1022 (0.1%)	0.93	2/1377 (0.1%)
1	B	0.93	0/1012	0.95	2/1365 (0.1%)
1	C	0.85	1/1022 (0.1%)	0.95	3/1377 (0.2%)
1	D	0.91	0/1012	1.00	2/1365 (0.1%)
1	E	0.80	0/1007	0.86	0/1358
1	F	0.87	0/1001	0.97	3/1350 (0.2%)
1	G	0.79	0/999	0.85	1/1347 (0.1%)
1	H	0.76	0/1001	0.83	0/1350
2	I	0.86	0/130	0.74	0/178
2	J	0.99	0/130	0.82	0/178
2	K	0.83	0/130	0.67	0/178
2	L	0.79	0/119	0.80	0/164
2	M	0.78	0/130	0.72	0/178
2	N	0.79	0/119	0.78	0/164
2	O	0.76	0/130	0.75	0/178
2	P	0.79	0/130	0.73	0/178
All	All	0.85	2/9094 (0.0%)	0.90	13/12285 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
1	D	0	1
1	E	0	1
2	I	0	1
2	J	0	1
All	All	0	5

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	53	CYS	CB-SG	-5.30	1.73	1.81
1	A	16	SER	CB-OG	-5.06	1.35	1.42

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	124	PHE	CB-CG-CD2	10.34	128.03	120.80
1	C	33	ASP	CB-CG-OD2	7.34	124.91	118.30
1	C	91	TYR	CB-CG-CD1	6.86	125.11	121.00
1	F	88	ASP	CB-CG-OD2	6.62	124.26	118.30
1	D	124	PHE	CB-CG-CD1	-6.01	116.59	120.80
1	C	91	TYR	CB-CG-CD2	-5.78	117.53	121.00
1	A	24	ARG	NE-CZ-NH1	-5.68	117.46	120.30
1	F	88	ASP	CB-CG-OD1	-5.44	113.40	118.30
1	B	35	VAL	CB-CA-C	-5.39	101.16	111.40
1	F	31	LEU	CB-CG-CD1	-5.35	101.90	111.00
1	B	51	MET	CG-SD-CE	5.14	108.42	100.20
1	G	63	ASP	CB-CG-OD2	-5.07	113.74	118.30
1	A	28	ARG	NE-CZ-NH2	-5.02	117.79	120.30

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	31	LEU	Mainchain
1	D	124	PHE	Sidechain
1	E	31	LEU	Mainchain
2	I	513	GLY	Peptide
2	J	513	GLY	Peptide

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	1008	0	1023	36	0
1	B	998	0	1017	37	0
1	C	1008	0	1023	40	0
1	D	998	0	1017	40	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	993	0	1012	35	0
1	F	987	0	1007	52	0
1	G	985	0	1008	24	0
1	H	987	0	1007	44	0
2	I	126	0	124	3	0
2	J	126	0	124	5	0
2	K	126	0	124	13	0
2	L	115	0	111	6	0
2	M	126	0	124	7	0
2	N	115	0	111	6	0
2	O	126	0	124	7	0
2	P	126	0	124	4	0
3	A	26	0	0	4	0
3	B	23	0	0	2	0
3	C	11	0	0	1	0
3	D	23	0	0	0	0
3	E	12	0	0	2	0
3	F	15	0	0	2	0
3	G	15	0	0	1	0
3	H	7	0	0	3	0
3	I	1	0	0	0	0
3	J	3	0	0	0	0
3	K	1	0	0	0	0
3	L	1	0	0	1	0
3	M	1	0	0	0	0
3	N	1	0	0	0	0
3	O	1	0	0	0	0
All	All	9091	0	9080	282	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 16.

All (282) 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:A:8:GLN:HG3	1:B:96:ASN:ND2	1.49	1.25
1:E:8:GLN:HG3	1:F:96:ASN:ND2	1.56	1.19
1:A:24:ARG:HG3	1:A:24:ARG:NH1	1.40	1.15
1:A:24:ARG:HH11	1:A:24:ARG:CG	1.57	1.13
1:H:43:PHE:HE2	1:H:84:ASN:HD22	0.98	0.94
1:B:110:MET:HG2	1:H:110:MET:CE	1.99	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:110:MET:HG2	1:H:110:MET:HE2	1.51	0.92
1:B:69:LEU:HD13	1:D:69:LEU:HD13	1.51	0.91
1:C:62:THR:CG2	1:E:28:ARG:HE	1.85	0.88
1:E:96:ASN:ND2	1:F:8:GLN:HG2	1.90	0.87
1:E:8:GLN:HG3	1:F:96:ASN:HD22	1.36	0.86
2:O:510:VAL:CG1	1:H:24:ARG:HH22	1.87	0.86
1:E:8:GLN:HG3	1:F:96:ASN:HD21	1.41	0.85
1:D:58:TYR:O	1:D:62:THR:HB	1.77	0.84
1:D:23:ASN:HD21	1:D:26:ARG:HH21	1.25	0.84
1:A:8:GLN:HG3	1:B:96:ASN:HD21	1.41	0.83
1:H:106:MET:HE2	1:H:121:CYS:HB3	1.60	0.83
1:B:69:LEU:CD1	1:D:69:LEU:HD13	2.09	0.82
1:A:58:TYR:O	1:A:62:THR:HB	1.79	0.82
1:G:96:ASN:ND2	1:H:8:GLN:HG3	1.96	0.81
1:D:126:LYS:HE3	1:D:126:LYS:HA	1.64	0.79
1:F:92:THR:O	1:F:93:SER:HB2	1.80	0.79
1:F:58:TYR:O	1:F:62:THR:HB	1.83	0.79
2:I:498:ARG:HB2	1:B:6:ASP:HB2	1.65	0.77
1:H:43:PHE:HE2	1:H:84:ASN:ND2	1.78	0.76
1:D:35:VAL:HG23	1:D:44:ARG:NH1	2.01	0.76
1:C:23:ASN:HD21	1:C:26:ARG:HH21	1.32	0.75
1:A:23:ASN:HD21	1:A:26:ARG:HH21	1.34	0.75
2:J:513:GLY:O	2:J:514:PRO:O	2.05	0.75
1:F:81:GLU:O	1:F:85:ILE:HG13	1.86	0.74
1:G:66:LYS:HA	1:G:69:LEU:HD22	1.68	0.74
2:P:512:PRO:C	2:P:514:PRO:HD3	2.07	0.74
1:G:31:LEU:HD12	1:H:51:MET:SD	2.28	0.73
3:G:130:HOH:O	1:H:95:LEU:HB2	1.87	0.73
1:B:122:ARG:NH1	3:B:151:HOH:O	2.22	0.72
1:C:13:ARG:HG2	3:L:74:HOH:O	1.89	0.72
1:A:24:ARG:HH11	1:A:24:ARG:HG3	0.66	0.71
2:K:498:ARG:HB2	1:D:6:ASP:HB3	1.72	0.71
1:C:58:TYR:O	1:C:62:THR:HB	1.91	0.71
2:O:510:VAL:CG1	1:H:24:ARG:NH2	2.56	0.69
1:D:23:ASN:ND2	1:D:26:ARG:HH21	1.90	0.69
1:D:106:MET:HE2	1:D:121:CYS:HB3	1.75	0.69
1:B:92:THR:O	1:B:93:SER:HB2	1.92	0.69
1:D:24:ARG:HH11	1:D:24:ARG:HG3	1.58	0.69
1:E:8:GLN:OE1	1:F:94:ARG:NH1	2.26	0.68
1:A:24:ARG:CG	1:A:24:ARG:NH1	2.28	0.68
1:C:97:LEU:HD21	1:C:105:VAL:HG11	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:43:PHE:CE2	1:H:84:ASN:ND2	2.58	0.68
1:H:98:ARG:HH11	1:H:98:ARG:HB3	1.59	0.67
1:F:106:MET:HE3	1:F:121:CYS:HB3	1.77	0.67
1:E:94:ARG:HB3	1:F:8:GLN:HE22	1.59	0.67
1:D:35:VAL:CG2	1:D:44:ARG:HH11	2.07	0.67
1:H:61:PHE:O	1:H:67:ARG:HG3	1.95	0.67
1:D:24:ARG:HG3	1:D:24:ARG:NH1	2.09	0.67
2:O:510:VAL:HG13	1:H:24:ARG:HH22	1.60	0.67
1:A:106:MET:HE2	1:A:121:CYS:HB3	1.77	0.66
1:A:24:ARG:HD3	3:A:140:HOH:O	1.96	0.66
1:G:24:ARG:HG3	2:P:511:VAL:HB	1.76	0.66
2:O:504:THR:O	2:O:504:THR:HG22	1.95	0.66
1:B:20:LEU:O	1:B:24:ARG:HG2	1.95	0.66
1:B:110:MET:HG2	1:H:110:MET:HE1	1.77	0.66
1:F:106:MET:CE	1:F:121:CYS:HB3	2.26	0.66
1:C:64:GLN:H	1:C:64:GLN:CD	1.99	0.65
1:C:62:THR:HG23	1:E:28:ARG:NE	2.10	0.65
1:F:75:ASP:O	1:F:78:ILE:HG12	1.97	0.65
1:B:122:ARG:C	1:B:124:PHE:H	1.99	0.65
1:B:122:ARG:O	1:B:124:PHE:N	2.29	0.65
1:C:62:THR:CG2	1:E:28:ARG:NE	2.59	0.64
1:F:28:ARG:HB2	1:F:30:ILE:HD12	1.80	0.64
1:B:58:TYR:O	1:B:62:THR:HB	1.97	0.64
1:G:31:LEU:CD1	1:H:51:MET:SD	2.86	0.64
1:C:62:THR:HG23	1:E:28:ARG:HE	1.64	0.63
1:D:35:VAL:CG2	1:D:44:ARG:NH1	2.62	0.63
1:E:44:ARG:NH1	3:E:130:HOH:O	2.32	0.63
1:D:97:LEU:HD13	1:D:102:ILE:CD1	2.28	0.62
2:I:512:PRO:HB2	2:I:514:PRO:HD3	1.80	0.62
1:D:23:ASN:HD21	1:D:26:ARG:NH2	1.95	0.62
2:M:498:ARG:HA	1:F:6:ASP:O	1.98	0.62
1:D:75:ASP:OD1	1:D:77:GLU:HB2	1.98	0.62
1:E:96:ASN:HD22	1:F:8:GLN:HG2	1.62	0.62
1:D:92:THR:O	1:D:93:SER:HB2	2.00	0.61
1:G:47:LYS:O	1:G:51:MET:HG3	2.00	0.61
1:G:21:ASN:HA	1:G:24:ARG:HG2	1.83	0.61
1:D:66:LYS:HA	1:D:69:LEU:HD22	1.81	0.61
1:H:98:ARG:HB3	1:H:98:ARG:NH1	2.15	0.61
1:F:28:ARG:HB2	1:F:30:ILE:CD1	2.29	0.61
1:G:64:GLN:OE1	1:G:64:GLN:N	2.32	0.61
1:H:35:VAL:HB	3:H:137:HOH:O	2.01	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:504:THR:O	2:K:504:THR:HG22	2.01	0.60
2:O:513:GLY:N	2:O:514:PRO:HD3	2.15	0.60
2:L:512:PRO:C	2:L:514:PRO:HD2	2.21	0.60
1:D:126:LYS:HA	1:D:126:LYS:CE	2.30	0.60
1:D:97:LEU:HD13	1:D:102:ILE:HD12	1.84	0.60
1:F:40:ARG:NH2	3:F:142:HOH:O	2.22	0.59
1:H:92:THR:O	1:H:93:SER:HB2	2.03	0.59
1:A:98:ARG:HG3	1:B:6:ASP:OD1	2.03	0.59
1:C:64:GLN:HE22	2:N:514:PRO:HD2	1.68	0.59
1:B:110:MET:CG	1:H:110:MET:HE2	2.28	0.59
1:G:53:CYS:SG	1:G:90:MET:HG2	2.43	0.58
1:G:51:MET:SD	1:H:31:LEU:HD12	2.44	0.58
1:G:47:LYS:NZ	1:G:67:ARG:O	2.37	0.58
1:C:92:THR:O	1:C:93:SER:HB2	2.04	0.58
1:E:96:ASN:ND2	1:F:8:GLN:HE21	2.02	0.57
1:E:92:THR:O	1:E:93:SER:HB2	2.04	0.57
1:A:41:GLU:HG3	1:A:80:PRO:HB3	1.87	0.57
2:N:513:GLY:O	2:N:514:PRO:C	2.41	0.57
1:A:62:THR:HG21	3:A:146:HOH:O	2.05	0.57
1:B:6:ASP:O	1:B:7:SER:HB2	2.05	0.56
1:B:69:LEU:HD13	1:D:69:LEU:CD1	2.29	0.56
1:C:64:GLN:NE2	2:N:514:PRO:HG2	2.21	0.56
1:A:51:MET:SD	1:B:31:LEU:HD12	2.46	0.56
1:D:127:ALA:O	1:D:128:SER:HB2	2.06	0.56
1:G:106:MET:HE3	1:G:125:ILE:HD11	1.88	0.56
2:L:513:GLY:N	2:L:514:PRO:HD2	2.21	0.56
1:G:58:TYR:O	1:G:62:THR:HB	2.06	0.56
1:B:126:LYS:HA	1:B:126:LYS:HE3	1.88	0.55
2:O:510:VAL:HG13	1:H:24:ARG:NH2	2.22	0.55
1:H:35:VAL:HG22	1:H:44:ARG:HD2	1.88	0.55
1:A:92:THR:O	1:A:93:SER:HB2	2.07	0.55
1:H:122:ARG:O	1:H:124:PHE:N	2.40	0.55
1:C:8:GLN:HB2	1:D:96:ASN:ND2	2.21	0.55
1:G:35:VAL:HG22	1:G:44:ARG:HH11	1.70	0.55
1:E:44:ARG:HD2	3:E:130:HOH:O	2.05	0.55
1:C:111:TYR:O	1:F:67:ARG:NH2	2.41	0.54
1:E:94:ARG:HB3	1:F:8:GLN:NE2	2.21	0.54
1:B:124:PHE:C	1:B:126:LYS:H	2.11	0.54
2:K:498:ARG:CB	1:D:6:ASP:HB3	2.37	0.54
2:P:512:PRO:HB2	2:P:514:PRO:HD3	1.88	0.54
2:K:513:GLY:N	2:K:514:PRO:CD	2.71	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:8:GLN:OE1	1:B:94:ARG:NH1	2.41	0.54
1:E:14:HIS:O	1:E:18:VAL:HG23	2.08	0.54
1:G:32:THR:HA	1:G:46:HIS:CD2	2.43	0.54
1:C:64:GLN:HE21	2:N:514:PRO:HG2	1.73	0.53
1:A:86:LEU:O	1:A:90:MET:HG3	2.09	0.53
1:G:93:SER:HA	1:H:14:HIS:CD2	2.44	0.53
2:M:513:GLY:N	2:M:514:PRO:CD	2.72	0.53
1:F:122:ARG:C	1:F:124:PHE:H	2.12	0.53
1:E:43:PHE:HE2	1:E:84:ASN:HD22	1.57	0.53
1:F:64:GLN:H	1:F:64:GLN:CD	2.12	0.52
1:C:96:ASN:ND2	1:D:8:GLN:HG3	2.24	0.52
2:J:512:PRO:HB2	2:J:514:PRO:HD3	1.90	0.52
1:F:123:LYS:O	1:F:123:LYS:HG3	2.08	0.52
1:A:66:LYS:HA	1:A:69:LEU:HD23	1.90	0.52
1:C:97:LEU:CD2	1:C:105:VAL:HG11	2.38	0.52
1:A:51:MET:SD	1:B:31:LEU:CD1	2.98	0.52
1:E:8:GLN:CG	1:F:96:ASN:ND2	2.50	0.52
1:A:62:THR:CG2	3:A:146:HOH:O	2.57	0.52
1:B:122:ARG:C	1:B:124:PHE:N	2.63	0.52
1:F:110:MET:HG2	1:F:111:TYR:N	2.24	0.52
1:G:28:ARG:HD2	2:P:513:GLY:HA2	1.92	0.51
1:H:62:THR:O	1:H:62:THR:HG22	2.09	0.51
1:C:46:HIS:CD2	1:D:48:THR:HG21	2.46	0.51
1:A:125:ILE:HG22	1:A:128:SER:OG	2.11	0.51
1:H:33:ASP:N	1:H:33:ASP:OD2	2.44	0.50
1:B:47:LYS:O	1:B:51:MET:HG3	2.11	0.50
1:E:98:ARG:HG2	1:F:6:ASP:OD1	2.12	0.50
1:B:6:ASP:O	1:B:7:SER:CB	2.59	0.50
2:M:504:THR:O	2:M:504:THR:HG22	2.12	0.50
1:F:35:VAL:HG13	1:F:44:ARG:HG2	1.93	0.49
1:D:127:ALA:O	1:D:128:SER:CB	2.60	0.49
1:G:126:LYS:HE3	1:G:126:LYS:HA	1.93	0.49
1:A:20:LEU:O	1:A:24:ARG:HG2	2.13	0.49
1:F:106:MET:HE2	1:F:121:CYS:C	2.33	0.49
2:K:513:GLY:N	2:K:514:PRO:HD3	2.27	0.49
1:C:103:MET:HA	1:C:103:MET:CE	2.42	0.49
1:C:122:ARG:C	1:C:124:PHE:H	2.16	0.49
1:E:106:MET:HE2	1:E:121:CYS:HB3	1.95	0.49
1:E:96:ASN:HD21	1:F:8:GLN:HG2	1.72	0.48
1:D:92:THR:O	1:D:93:SER:CB	2.61	0.48
2:O:504:THR:O	2:O:504:THR:CG2	2.61	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:126:LYS:C	1:B:128:SER:H	2.17	0.48
1:A:96:ASN:ND2	1:B:8:GLN:HG2	2.28	0.48
1:E:125:ILE:HG22	1:E:125:ILE:O	2.13	0.48
2:K:501:ILE:HG13	2:K:502:ILE:N	2.27	0.48
2:M:512:PRO:HB2	1:F:28:ARG:HH12	1.78	0.48
1:A:94:ARG:HB3	1:B:8:GLN:NE2	2.28	0.48
1:E:8:GLN:CG	1:F:96:ASN:HD22	2.17	0.47
1:C:13:ARG:HD3	2:L:504:THR:O	2.13	0.47
1:C:64:GLN:HE22	2:N:514:PRO:CD	2.27	0.47
1:C:75:ASP:HB3	1:C:78:ILE:HG12	1.96	0.47
1:E:6:ASP:HA	1:F:98:ARG:HB3	1.95	0.47
1:F:120:THR:O	1:F:124:PHE:HB2	2.15	0.47
1:E:6:ASP:OD1	1:F:98:ARG:HD3	2.14	0.47
1:F:43:PHE:HE2	1:F:84:ASN:HD22	1.63	0.46
1:H:44:ARG:O	3:H:133:HOH:O	2.20	0.46
1:E:21:ASN:O	1:E:25:LEU:HG	2.16	0.46
1:F:106:MET:CE	1:F:121:CYS:CB	2.93	0.46
1:C:21:ASN:HA	1:C:24:ARG:HG2	1.96	0.46
1:A:38:VAL:HG22	1:A:74:LEU:HD12	1.96	0.46
1:C:124:PHE:C	1:C:126:LYS:H	2.18	0.46
1:E:124:PHE:C	1:E:126:LYS:H	2.19	0.46
1:C:35:VAL:HG22	1:C:44:ARG:HD2	1.97	0.46
1:F:20:LEU:O	1:F:24:ARG:HG2	2.16	0.46
1:F:123:LYS:O	1:F:123:LYS:CG	2.64	0.46
1:C:118:VAL:O	1:C:121:CYS:HB2	2.16	0.45
1:D:20:LEU:O	1:D:24:ARG:HG2	2.16	0.45
1:F:106:MET:HE2	1:F:121:CYS:CB	2.47	0.45
1:E:106:MET:HB2	1:E:106:MET:HE2	1.60	0.45
1:F:106:MET:HE2	1:F:121:CYS:HB3	1.99	0.45
1:C:86:LEU:HD12	1:C:108:THR:HB	1.99	0.45
1:A:23:ASN:ND2	1:A:26:ARG:HE	2.15	0.45
2:I:504:THR:HG22	2:I:504:THR:O	2.16	0.45
2:M:501:ILE:HG13	2:M:502:ILE:N	2.31	0.45
1:H:79:ASN:OD1	1:H:104:ALA:HB1	2.17	0.45
1:B:110:MET:HE1	1:H:122:ARG:NE	2.32	0.45
1:H:11:PHE:HB2	1:H:14:HIS:HB2	1.99	0.45
1:B:88:ASP:OD2	3:B:141:HOH:O	2.21	0.44
2:K:498:ARG:CA	1:D:6:ASP:HB3	2.47	0.44
1:A:123:LYS:O	1:A:123:LYS:CG	2.64	0.44
2:K:510:VAL:CG1	1:D:24:ARG:HH12	2.30	0.44
1:B:106:MET:HE2	1:B:121:CYS:HB3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:67:ARG:NH1	3:F:134:HOH:O	2.49	0.44
1:F:122:ARG:O	1:F:124:PHE:N	2.51	0.44
1:C:106:MET:CE	1:C:121:CYS:HB3	2.48	0.44
1:H:47:LYS:O	1:H:51:MET:HG3	2.18	0.44
1:A:6:ASP:OD1	1:B:98:ARG:HD2	2.17	0.44
3:C:137:HOH:O	1:F:62:THR:CG2	2.65	0.44
1:H:62:THR:CG2	1:H:62:THR:O	2.65	0.44
1:F:85:ILE:HD13	1:F:96:ASN:O	2.17	0.43
1:A:125:ILE:CG2	1:A:128:SER:OG	2.66	0.43
1:D:24:ARG:HH11	1:D:24:ARG:CG	2.24	0.43
1:D:63:ASP:OD1	1:D:63:ASP:C	2.57	0.43
1:F:23:ASN:HD21	1:F:26:ARG:HH21	1.66	0.43
1:F:28:ARG:CB	1:F:30:ILE:HD12	2.48	0.43
1:G:92:THR:O	1:G:93:SER:CB	2.67	0.43
1:C:96:ASN:HD21	1:D:8:GLN:HG3	1.83	0.43
2:L:513:GLY:N	2:L:514:PRO:CD	2.82	0.43
1:C:8:GLN:CD	1:D:94:ARG:HG2	2.39	0.43
1:H:122:ARG:C	1:H:124:PHE:N	2.72	0.43
2:M:512:PRO:CB	2:M:514:PRO:HD2	2.49	0.42
1:E:121:CYS:O	1:E:124:PHE:HB3	2.20	0.42
2:J:512:PRO:HB2	2:J:514:PRO:CD	2.49	0.42
1:F:75:ASP:HA	1:F:76:PRO:HD3	1.80	0.42
1:E:53:CYS:SG	1:E:90:MET:HG2	2.59	0.42
1:G:96:ASN:CG	1:H:8:GLN:HG3	2.39	0.42
1:B:6:ASP:HB3	1:B:7:SER:H	1.68	0.42
1:A:126:LYS:HA	1:A:126:LYS:HE3	2.00	0.42
1:H:23:ASN:HD21	1:H:26:ARG:HH21	1.68	0.42
1:H:106:MET:HE2	1:H:106:MET:HB2	1.89	0.42
2:K:513:GLY:HA2	1:D:28:ARG:HD2	2.02	0.42
2:K:500:GLU:H	2:K:500:GLU:HG3	1.67	0.42
1:H:35:VAL:CG2	3:H:137:HOH:O	2.68	0.42
1:C:106:MET:HB2	1:C:106:MET:HE2	1.88	0.42
1:G:35:VAL:HG22	1:G:44:ARG:NH1	2.34	0.42
1:B:110:MET:HE2	1:H:122:ARG:HE	1.85	0.42
1:E:87:LEU:HD12	1:E:87:LEU:HA	1.85	0.42
1:E:66:LYS:HA	1:E:69:LEU:HD22	2.02	0.42
2:M:512:PRO:C	2:M:514:PRO:HD2	2.40	0.41
1:G:75:ASP:HA	1:G:76:PRO:HD3	1.79	0.41
2:K:504:THR:O	2:K:504:THR:CG2	2.68	0.41
1:H:122:ARG:O	1:H:123:LYS:C	2.59	0.41
2:K:498:ARG:HA	1:D:6:ASP:C	2.41	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:ILE:HG12	2:J:501:ILE:HG12	2.03	0.41
1:B:110:MET:CE	1:H:122:ARG:NE	2.84	0.41
1:A:23:ASN:HD22	1:A:26:ARG:HE	1.67	0.41
1:C:86:LEU:HD13	1:C:112:LEU:HD12	2.03	0.41
1:A:75:ASP:HA	1:A:76:PRO:HD3	1.84	0.41
1:A:26:ARG:HD3	1:A:91:TYR:CE1	2.55	0.41
1:C:13:ARG:CD	2:L:504:THR:O	2.69	0.41
1:H:110:MET:HE3	1:H:110:MET:HB2	1.96	0.41
1:F:28:ARG:CB	1:F:30:ILE:CD1	2.98	0.41
1:A:124:PHE:O	1:A:126:LYS:N	2.54	0.41
1:C:106:MET:HE2	1:C:121:CYS:HB3	2.03	0.41
1:E:11:PHE:CD1	1:F:89:PHE:HE1	2.38	0.41
1:H:20:LEU:HG	1:H:20:LEU:O	2.17	0.41
1:B:79:ASN:HA	1:B:80:PRO:HD3	1.98	0.41
1:H:38:VAL:HG22	1:H:74:LEU:HD12	2.02	0.41
2:J:513:GLY:N	2:J:514:PRO:CD	2.84	0.40
1:C:6:ASP:OD2	1:D:98:ARG:HG3	2.21	0.40
1:F:38:VAL:HG13	1:F:78:ILE:HB	2.02	0.40
1:G:124:PHE:C	1:G:126:LYS:H	2.25	0.40
1:G:99:GLU:O	1:G:102:ILE:HG22	2.21	0.40
1:D:66:LYS:O	1:D:67:ARG:C	2.60	0.40
2:K:498:ARG:HA	1:D:6:ASP:O	2.21	0.40
1:C:28:ARG:NH1	2:L:512:PRO:O	2.52	0.40
1:C:55:GLY:HA3	1:E:29:ASP:OD1	2.21	0.40
1:F:47:LYS:O	1:F:51:MET:HG3	2.22	0.40
1:A:48:THR:HG23	3:A:153:HOH:O	2.21	0.40
2:N:513:GLY:N	2:N:514:PRO:HD3	2.37	0.40
1:C:103:MET:HE2	1:C:103:MET:HA	2.03	0.40
1:C:98:ARG:O	1:C:102:ILE:HB	2.22	0.40
1:F:38:VAL:CG1	1:F:78:ILE:HB	2.51	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	123/127 (97%)	111 (90%)	11 (9%)	1 (1%)	24	46
1	B	122/127 (96%)	112 (92%)	6 (5%)	4 (3%)	5	7
1	C	123/127 (97%)	112 (91%)	8 (6%)	3 (2%)	7	13
1	D	122/127 (96%)	109 (89%)	12 (10%)	1 (1%)	24	46
1	E	121/127 (95%)	113 (93%)	7 (6%)	1 (1%)	24	46
1	F	120/127 (94%)	111 (92%)	7 (6%)	2 (2%)	11	22
1	G	120/127 (94%)	110 (92%)	9 (8%)	1 (1%)	24	46
1	H	120/127 (94%)	108 (90%)	11 (9%)	1 (1%)	24	46
2	I	15/19 (79%)	15 (100%)	0	0	100	100
2	J	15/19 (79%)	15 (100%)	0	0	100	100
2	K	15/19 (79%)	14 (93%)	1 (7%)	0	100	100
2	L	14/19 (74%)	14 (100%)	0	0	100	100
2	M	15/19 (79%)	14 (93%)	1 (7%)	0	100	100
2	N	14/19 (74%)	14 (100%)	0	0	100	100
2	O	15/19 (79%)	13 (87%)	2 (13%)	0	100	100
2	P	15/19 (79%)	14 (93%)	1 (7%)	0	100	100
All	All	1089/1168 (93%)	999 (92%)	76 (7%)	14 (1%)	15	30

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	7	SER
1	B	123	LYS
1	A	125	ILE
1	C	123	LYS
1	C	125	ILE
1	F	123	LYS
1	G	125	ILE
1	D	126	LYS
1	E	125	ILE
1	H	123	LYS
1	B	125	ILE
1	B	127	ALA
1	F	39	SER
1	C	102	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	114/115 (99%)	97 (85%)	17 (15%)	4	6
1	B	113/115 (98%)	101 (89%)	12 (11%)	8	15
1	C	114/115 (99%)	97 (85%)	17 (15%)	4	6
1	D	113/115 (98%)	104 (92%)	9 (8%)	15	29
1	E	113/115 (98%)	97 (86%)	16 (14%)	4	7
1	F	112/115 (97%)	99 (88%)	13 (12%)	7	12
1	G	112/115 (97%)	101 (90%)	11 (10%)	10	19
1	H	112/115 (97%)	102 (91%)	10 (9%)	12	23
2	I	15/16 (94%)	14 (93%)	1 (7%)	20	40
2	J	15/16 (94%)	9 (60%)	6 (40%)	0	0
2	K	15/16 (94%)	14 (93%)	1 (7%)	20	40
2	L	14/16 (88%)	11 (79%)	3 (21%)	1	2
2	M	15/16 (94%)	14 (93%)	1 (7%)	20	40
2	N	14/16 (88%)	13 (93%)	1 (7%)	18	36
2	O	15/16 (94%)	11 (73%)	4 (27%)	0	1
2	P	15/16 (94%)	11 (73%)	4 (27%)	0	1
All	All	1021/1048 (97%)	895 (88%)	126 (12%)	6	11

All (126) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	GLN
1	A	10	GLN
1	A	13	ARG
1	A	20	LEU
1	A	24	ARG
1	A	31	LEU
1	A	54	SER
1	A	59	SER

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Mol	Chain	Res	Type
1	A	62	THR
1	A	67	ARG
1	A	94	ARG
1	A	98	ARG
1	A	102	ILE
1	A	115	GLU
1	A	125	ILE
1	A	126	LYS
1	A	129	GLU
2	I	500	GLU
1	B	7	SER
1	B	10	GLN
1	B	13	ARG
1	B	20	LEU
1	B	24	ARG
1	B	35	VAL
1	B	62	THR
1	B	64	GLN
1	B	69	LEU
1	B	98	ARG
1	B	106	MET
1	B	126	LYS
2	J	498	ARG
2	J	499	SER
2	J	500	GLU
2	J	501	ILE
2	J	502	ILE
2	J	504	THR
1	C	23	ASN
1	C	24	ARG
1	C	31	LEU
1	C	54	SER
1	C	62	THR
1	C	64	GLN
1	C	65	LEU
1	C	69	LEU
1	C	94	ARG
1	C	98	ARG
1	C	102	ILE
1	C	106	MET
1	C	110	MET
1	C	115	GLU

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Mol	Chain	Res	Type
1	C	119	ASP
1	C	123	LYS
1	C	129	GLU
2	K	500	GLU
1	D	8	GLN
1	D	31	LEU
1	D	62	THR
1	D	64	GLN
1	D	69	LEU
1	D	77	GLU
1	D	94	ARG
1	D	102	ILE
1	D	126	LYS
2	L	499	SER
2	L	500	GLU
2	L	504	THR
1	E	6	ASP
1	E	8	GLN
1	E	24	ARG
1	E	31	LEU
1	E	35	VAL
1	E	42	GLN
1	E	54	SER
1	E	64	GLN
1	E	67	ARG
1	E	69	LEU
1	E	70	SER
1	E	86	LEU
1	E	106	MET
1	E	115	GLU
1	E	126	LYS
1	E	128	SER
2	M	500	GLU
1	F	8	GLN
1	F	13	ARG
1	F	20	LEU
1	F	35	VAL
1	F	38	VAL
1	F	41	GLU
1	F	54	SER
1	F	62	THR
1	F	64	GLN

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Mol	Chain	Res	Type
1	F	93	SER
1	F	102	ILE
1	F	110	MET
1	F	126	LYS
2	N	499	SER
1	G	13	ARG
1	G	24	ARG
1	G	27	SER
1	G	40	ARG
1	G	54	SER
1	G	62	THR
1	G	64	GLN
1	G	69	LEU
1	G	86	LEU
1	G	125	ILE
1	G	126	LYS
2	O	498	ARG
2	O	499	SER
2	O	500	GLU
2	O	501	ILE
1	H	6	ASP
1	H	8	GLN
1	H	13	ARG
1	H	24	ARG
1	H	54	SER
1	H	64	GLN
1	H	65	LEU
1	H	103	MET
1	H	125	ILE
1	H	126	LYS
2	P	498	ARG
2	P	499	SER
2	P	504	THR
2	P	506	PRO

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	ASN
1	A	96	ASN
1	B	8	GLN
1	B	10	GLN

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Mol	Chain	Res	Type
1	B	23	ASN
1	B	96	ASN
1	C	10	GLN
1	C	23	ASN
1	C	64	GLN
1	C	96	ASN
1	C	113	GLN
1	D	23	ASN
1	D	96	ASN
1	E	23	ASN
1	E	84	ASN
1	E	96	ASN
1	F	8	GLN
1	F	23	ASN
1	F	84	ASN
1	F	96	ASN
1	F	101	ASN
1	G	8	GLN
1	G	23	ASN
1	G	84	ASN
1	G	96	ASN
1	H	23	ASN
1	H	96	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 carbohydrates 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	125/127 (98%)	-0.09	4 (3%)	51	44	40, 58, 98, 124	0
1	B	124/127 (97%)	-0.04	3 (2%)	62	56	41, 56, 95, 116	0
1	C	125/127 (98%)	0.30	8 (6%)	23	17	46, 70, 112, 134	0
1	D	124/127 (97%)	-0.16	4 (3%)	51	44	43, 55, 98, 114	0
1	E	123/127 (96%)	0.06	3 (2%)	62	56	48, 67, 97, 125	0
1	F	122/127 (96%)	0.02	6 (4%)	33	26	50, 67, 100, 115	0
1	G	122/127 (96%)	-0.06	6 (4%)	33	26	50, 72, 103, 125	0
1	H	122/127 (96%)	0.03	5 (4%)	41	33	55, 75, 107, 117	0
2	I	17/19 (89%)	1.27	3 (17%)	2	1	59, 70, 109, 116	0
2	J	17/19 (89%)	0.31	3 (17%)	2	1	49, 58, 105, 113	0
2	K	17/19 (89%)	1.25	5 (29%)	1	0	61, 81, 119, 122	0
2	L	16/19 (84%)	0.42	3 (18%)	2	1	64, 77, 104, 104	0
2	M	17/19 (89%)	1.14	4 (23%)	1	0	68, 82, 115, 122	0
2	N	16/19 (84%)	0.99	2 (12%)	5	3	61, 78, 107, 107	0
2	O	17/19 (89%)	1.50	6 (35%)	0	0	68, 93, 123, 126	0
2	P	17/19 (89%)	1.09	5 (29%)	1	0	79, 97, 122, 127	0
All	All	1121/1168 (95%)	0.13	70 (6%)	24	18	40, 68, 109, 134	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	I	499	SER	8.4
2	K	499	SER	7.1
2	P	514	PRO	6.8
2	M	498	ARG	6.5
2	O	499	SER	6.2

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Mol	Chain	Res	Type	RSRZ
2	J	498	ARG	6.0
1	D	5	ALA	5.9
2	M	499	SER	5.8
2	O	498	ARG	5.8
2	O	502	ILE	5.0
2	M	514	PRO	4.8
1	F	5	ALA	4.8
2	O	500	GLU	4.6
1	A	5	ALA	4.4
2	P	499	SER	4.4
2	N	499	SER	4.3
1	A	127	ALA	4.2
2	N	500	GLU	4.2
1	C	128	SER	4.0
1	D	6	ASP	4.0
1	C	129	GLU	4.0
1	C	98	ARG	3.8
2	L	499	SER	3.8
2	P	498	ARG	3.7
1	C	40	ARG	3.6
2	O	501	ILE	3.6
2	K	498	ARG	3.4
2	K	514	PRO	3.4
2	I	498	ARG	3.4
2	L	514	PRO	3.3
2	K	500	GLU	3.3
1	C	97	LEU	3.2
1	F	98	ARG	3.1
1	H	7	SER	3.1
2	I	500	GLU	3.1
1	D	7	SER	3.0
1	E	6	ASP	3.0
1	C	127	ALA	2.9
1	F	124	PHE	2.8
1	H	98	ARG	2.8
1	H	6	ASP	2.8
1	A	128	SER	2.8
1	H	64	GLN	2.8
1	C	126	LYS	2.7
1	G	128	SER	2.7
1	F	6	ASP	2.7
2	J	499	SER	2.7

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Mol	Chain	Res	Type	RSRZ
1	E	126	LYS	2.7
2	O	514	PRO	2.7
1	B	5	ALA	2.7
1	C	99	GLU	2.7
1	D	127	ALA	2.6
1	G	99	GLU	2.6
2	K	501	ILE	2.6
1	E	128	SER	2.6
1	A	6	ASP	2.6
1	G	98	ARG	2.4
1	B	6	ASP	2.4
2	P	500	GLU	2.4
1	G	127	ALA	2.4
2	L	513	GLY	2.3
2	M	502	ILE	2.3
1	F	101	ASN	2.3
2	J	500	GLU	2.2
1	H	8	GLN	2.1
2	P	513	GLY	2.1
1	F	99	GLU	2.1
1	B	128	SER	2.1
1	G	124	PHE	2.0
1	G	40	ARG	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.