



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:54 PM GMT

PDB ID : 4N0O
Title : Complex structure of Arterivirus nonstructural protein 10 (helicase) with DNA
Authors : Deng, Z.; Chen, Z.
Deposited on : 2013-10-02
Resolution : 2.65 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:

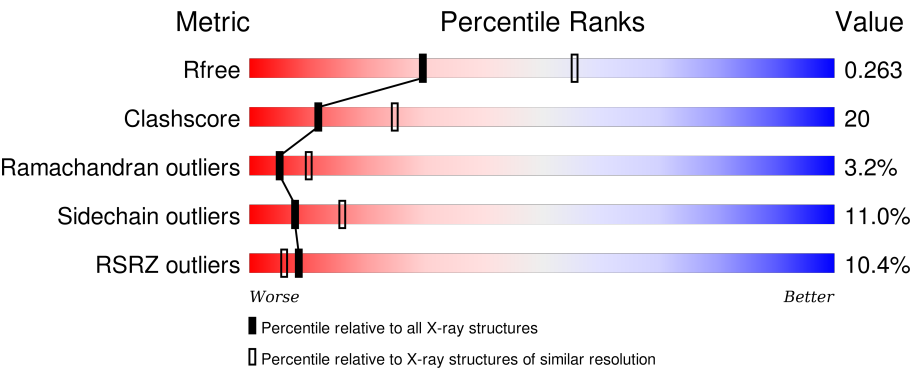
MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : 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 2.65 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.



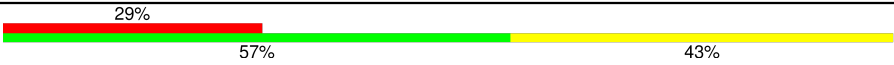


Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	91344	3152 (2.70-2.62)
Clashscore	102246	3524 (2.70-2.62)
Ramachandran outliers	100387	3469 (2.70-2.62)
Sidechain outliers	100360	3469 (2.70-2.62)
RSRZ outliers	91569	3161 (2.70-2.62)

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	423	
1	C	423	
1	E	423	
1	G	423	
2	B	7	

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Mol	Chain	Length	Quality of chain
2	D	7	
2	F	7	
2	H	7	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	ZN	G	503	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 12998 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 Replicase polypeptide 1ab.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	395	Total	C	N	O	S	0	0	0
			2897	1850	508	519	20			
1	C	398	Total	C	N	O	S	0	0	0
			2916	1867	506	524	19			
1	E	396	Total	C	N	O	S	0	0	0
			2896	1850	505	523	18			
1	G	393	Total	C	N	O	S	0	0	0
			2889	1845	509	517	18			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	EXPRESSION TAG	UNP P19811
A	-19	GLY	-	EXPRESSION TAG	UNP P19811
A	-18	SER	-	EXPRESSION TAG	UNP P19811
A	-17	SER	-	EXPRESSION TAG	UNP P19811
A	-16	HIS	-	EXPRESSION TAG	UNP P19811
A	-15	HIS	-	EXPRESSION TAG	UNP P19811
A	-14	HIS	-	EXPRESSION TAG	UNP P19811
A	-13	HIS	-	EXPRESSION TAG	UNP P19811
A	-12	HIS	-	EXPRESSION TAG	UNP P19811
A	-11	SER	-	EXPRESSION TAG	UNP P19811
A	-10	SER	-	EXPRESSION TAG	UNP P19811
A	-9	GLY	-	EXPRESSION TAG	UNP P19811
A	-8	GLU	-	EXPRESSION TAG	UNP P19811
A	-7	ASN	-	EXPRESSION TAG	UNP P19811
A	-6	LEU	-	EXPRESSION TAG	UNP P19811
A	-5	TYR	-	EXPRESSION TAG	UNP P19811
A	-4	PHE	-	EXPRESSION TAG	UNP P19811
A	-3	GLN	-	EXPRESSION TAG	UNP P19811
A	-2	GLY	-	EXPRESSION TAG	UNP P19811
A	-1	HIS	-	EXPRESSION TAG	UNP P19811
A	0	MET	-	EXPRESSION TAG	UNP P19811

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-20	MET	-	EXPRESSION TAG	UNP P19811
C	-19	GLY	-	EXPRESSION TAG	UNP P19811
C	-18	SER	-	EXPRESSION TAG	UNP P19811
C	-17	SER	-	EXPRESSION TAG	UNP P19811
C	-16	HIS	-	EXPRESSION TAG	UNP P19811
C	-15	HIS	-	EXPRESSION TAG	UNP P19811
C	-14	HIS	-	EXPRESSION TAG	UNP P19811
C	-13	HIS	-	EXPRESSION TAG	UNP P19811
C	-12	HIS	-	EXPRESSION TAG	UNP P19811
C	-11	SER	-	EXPRESSION TAG	UNP P19811
C	-10	SER	-	EXPRESSION TAG	UNP P19811
C	-9	GLY	-	EXPRESSION TAG	UNP P19811
C	-8	GLU	-	EXPRESSION TAG	UNP P19811
C	-7	ASN	-	EXPRESSION TAG	UNP P19811
C	-6	LEU	-	EXPRESSION TAG	UNP P19811
C	-5	TYR	-	EXPRESSION TAG	UNP P19811
C	-4	PHE	-	EXPRESSION TAG	UNP P19811
C	-3	GLN	-	EXPRESSION TAG	UNP P19811
C	-2	GLY	-	EXPRESSION TAG	UNP P19811
C	-1	HIS	-	EXPRESSION TAG	UNP P19811
C	0	MET	-	EXPRESSION TAG	UNP P19811
E	-20	MET	-	EXPRESSION TAG	UNP P19811
E	-19	GLY	-	EXPRESSION TAG	UNP P19811
E	-18	SER	-	EXPRESSION TAG	UNP P19811
E	-17	SER	-	EXPRESSION TAG	UNP P19811
E	-16	HIS	-	EXPRESSION TAG	UNP P19811
E	-15	HIS	-	EXPRESSION TAG	UNP P19811
E	-14	HIS	-	EXPRESSION TAG	UNP P19811
E	-13	HIS	-	EXPRESSION TAG	UNP P19811
E	-12	HIS	-	EXPRESSION TAG	UNP P19811
E	-11	SER	-	EXPRESSION TAG	UNP P19811
E	-10	SER	-	EXPRESSION TAG	UNP P19811
E	-9	GLY	-	EXPRESSION TAG	UNP P19811
E	-8	GLU	-	EXPRESSION TAG	UNP P19811
E	-7	ASN	-	EXPRESSION TAG	UNP P19811
E	-6	LEU	-	EXPRESSION TAG	UNP P19811
E	-5	TYR	-	EXPRESSION TAG	UNP P19811
E	-4	PHE	-	EXPRESSION TAG	UNP P19811
E	-3	GLN	-	EXPRESSION TAG	UNP P19811
E	-2	GLY	-	EXPRESSION TAG	UNP P19811
E	-1	HIS	-	EXPRESSION TAG	UNP P19811
E	0	MET	-	EXPRESSION TAG	UNP P19811

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-20	MET	-	EXPRESSION TAG	UNP P19811
G	-19	GLY	-	EXPRESSION TAG	UNP P19811
G	-18	SER	-	EXPRESSION TAG	UNP P19811
G	-17	SER	-	EXPRESSION TAG	UNP P19811
G	-16	HIS	-	EXPRESSION TAG	UNP P19811
G	-15	HIS	-	EXPRESSION TAG	UNP P19811
G	-14	HIS	-	EXPRESSION TAG	UNP P19811
G	-13	HIS	-	EXPRESSION TAG	UNP P19811
G	-12	HIS	-	EXPRESSION TAG	UNP P19811
G	-11	SER	-	EXPRESSION TAG	UNP P19811
G	-10	SER	-	EXPRESSION TAG	UNP P19811
G	-9	GLY	-	EXPRESSION TAG	UNP P19811
G	-8	GLU	-	EXPRESSION TAG	UNP P19811
G	-7	ASN	-	EXPRESSION TAG	UNP P19811
G	-6	LEU	-	EXPRESSION TAG	UNP P19811
G	-5	TYR	-	EXPRESSION TAG	UNP P19811
G	-4	PHE	-	EXPRESSION TAG	UNP P19811
G	-3	GLN	-	EXPRESSION TAG	UNP P19811
G	-2	GLY	-	EXPRESSION TAG	UNP P19811
G	-1	HIS	-	EXPRESSION TAG	UNP P19811
G	0	MET	-	EXPRESSION TAG	UNP P19811

- Molecule 2 is a DNA chain called DNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	7	Total 137	C 70	N 14	O 47	P 6	0	0	0
2	D	7	Total 137	C 70	N 14	O 47	P 6	0	0	0
2	F	7	Total 137	C 70	N 14	O 47	P 6	0	0	0
2	H	7	Total 137	C 70	N 14	O 47	P 6	0	0	0

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	3	Total	Zn	0	0
			3	3		
3	A	3	Total	Zn	0	0
			3	3		
3	C	3	Total	Zn	0	0
			3	3		

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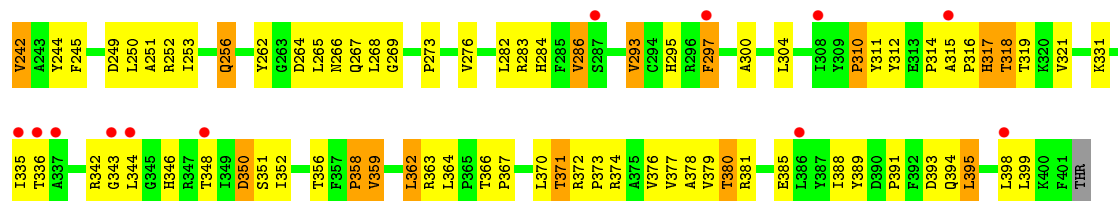
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	3	Total 3	Zn 3	0	0

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

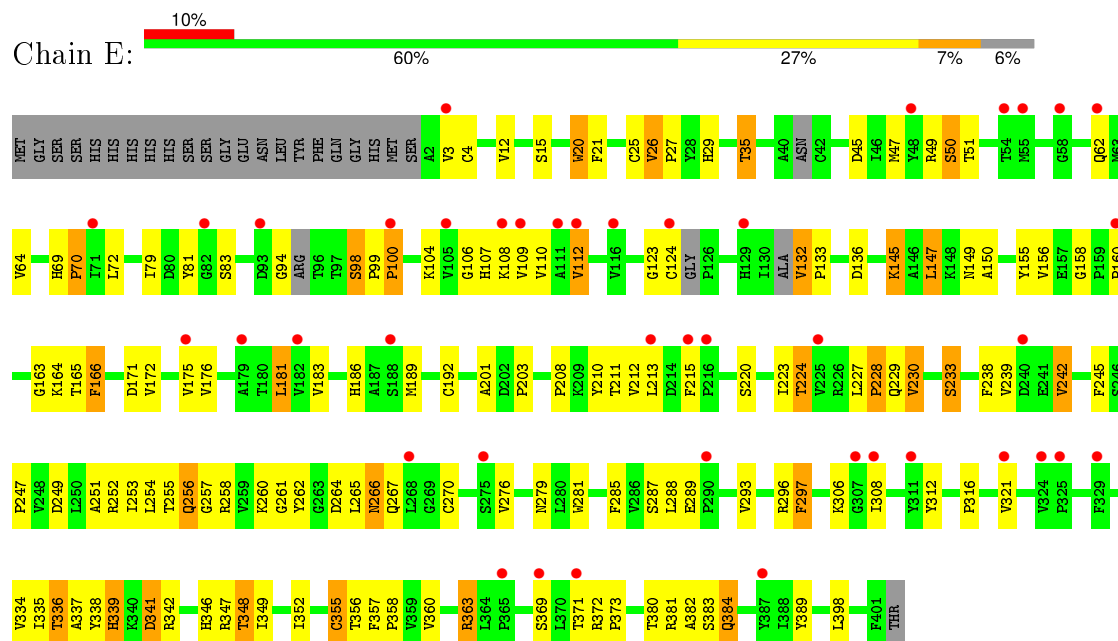
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	G	1	Total 1	Ca 1	0	0
4	A	2	Total 2	Ca 2	0	0
4	C	2	Total 2	Ca 2	0	0
4	E	2	Total 2	Ca 2	0	0

- Molecule 5 is water.

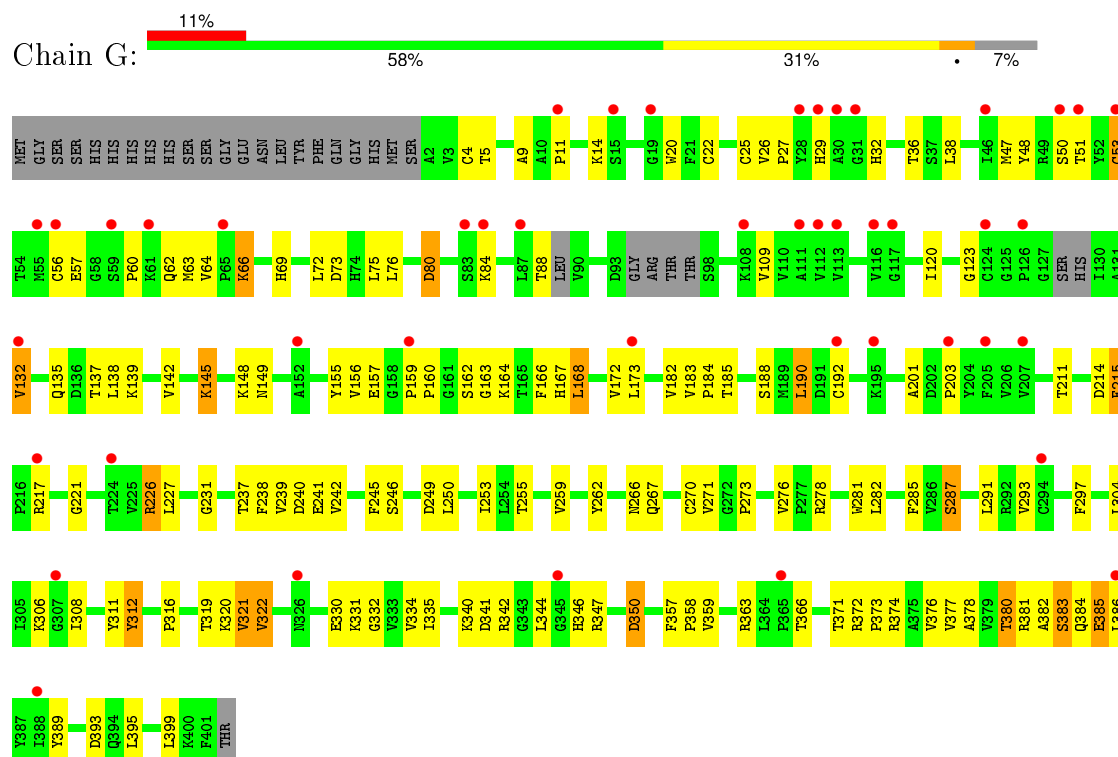
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	234	Total 234	O 234	0	0
5	C	169	Total 169	O 169	0	0
5	E	186	Total 186	O 186	0	0
5	G	216	Total 216	O 216	0	0
5	B	6	Total 6	O 6	0	0
5	D	4	Total 4	O 4	0	0
5	F	12	Total 12	O 12	0	0
5	H	6	Total 6	O 6	0	0



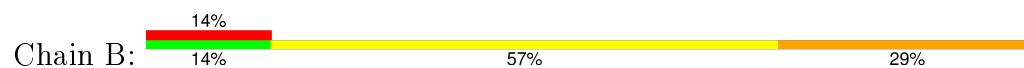
● Molecule 1: Replicase polypeptide 1ab



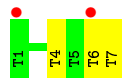
● Molecule 1: Replicase polypeptide 1ab



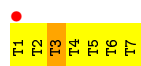
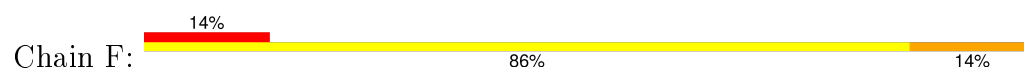
● Molecule 2: DNA



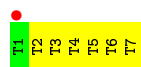
● Molecule 2: DNA



● Molecule 2: DNA



● Molecule 2: DNA



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	56.63Å 88.84Å 128.78Å 81.74° 90.01° 71.42°	Depositor
Resolution (Å)	50.00 – 2.65 42.43 – 2.63	Depositor EDS
% Data completeness (in resolution range)	95.7 (50.00-2.65) 90.6 (42.43-2.63)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.56 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.246 , 0.264 0.246 , 0.263	Depositor DCC
R_{free} test set	3305 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	54.0	Xtriage
Anisotropy	0.322	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 38.8	EDS
Estimated twinning fraction	0.532 for H, K, L 0.468 for H, H-K, -L 0.457 for h,h-k,-l	Xtriage
Reported twinning fraction	0.532 for H, K, L 0.468 for H, H-K, -L	Depositor
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 66828 reflections	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	12998	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 22.68 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 5.4307e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.51	2/2969 (0.1%)	0.70	2/4052 (0.0%)
1	C	0.50	1/2993 (0.0%)	0.65	0/4092
1	E	0.49	1/2970 (0.0%)	0.63	0/4057
1	G	0.49	1/2964 (0.0%)	0.64	1/4051 (0.0%)
2	B	0.31	0/150	1.04	2/230 (0.9%)
2	D	0.26	0/150	0.76	0/230
2	F	0.40	0/150	0.85	1/230 (0.4%)
2	H	0.32	0/150	0.79	0/230
All	All	0.49	5/12496 (0.0%)	0.67	6/17172 (0.0%)

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	C	0	1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	20	TRP	CD2-CE2	5.39	1.47	1.41
1	E	20	TRP	CD2-CE2	5.37	1.47	1.41
1	G	281	TRP	CD2-CE2	5.35	1.47	1.41
1	A	281	TRP	CD2-CE2	5.26	1.47	1.41
1	A	20	TRP	CD2-CE2	5.00	1.47	1.41

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	DT	P-O3'-C3'	9.78	131.44	119.70
2	B	4	DT	P-O3'-C3'	6.14	127.07	119.70
1	A	134	LEU	CA-CB-CG	5.37	127.66	115.30
1	G	190	LEU	CA-CB-CG	5.33	127.55	115.30
1	A	304	LEU	CA-CB-CG	5.27	127.43	115.30
2	F	3	DT	C1'-O4'-C4'	-5.14	104.96	110.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	1	SER	Peptide

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	2897	0	2802	129	0
1	C	2916	0	2810	116	0
1	E	2896	0	2790	123	0
1	G	2889	0	2811	105	0
2	B	137	0	86	6	0
2	D	137	0	86	3	0
2	F	137	0	86	9	0
2	H	137	0	86	9	0
3	A	3	0	0	0	0
3	C	3	0	0	0	0
3	E	3	0	0	0	0
3	G	3	0	0	2	0
4	A	2	0	0	0	0
4	C	2	0	0	0	0
4	E	2	0	0	0	0
4	G	1	0	0	0	0
5	A	234	0	0	3	0
5	B	6	0	0	0	0
5	C	169	0	0	4	0
5	D	4	0	0	0	0
5	E	186	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	F	12	0	0	1	0
5	G	216	0	0	8	0
5	H	6	0	0	0	0
All	All	12998	0	11557	480	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 20.

All (480) 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:1:DT:H2'	2:F:2:DT:H5'	1.15	1.12
2:H:4:DT:H2''	2:H:5:DT:H72	1.22	1.09
1:A:239:VAL:HG12	1:A:242:VAL:HG12	1.44	1.00
1:A:326:ASN:ND2	1:A:329:PHE:HB2	1.81	0.95
2:H:4:DT:H2''	2:H:5:DT:C7	1.96	0.95
1:E:172:VAL:HG21	1:E:181:LEU:HD13	1.49	0.93
1:E:358:PRO:HA	1:E:383:SER:HB2	1.52	0.91
1:E:242:VAL:HG11	1:E:261:GLY:HA3	1.52	0.90
2:F:1:DT:C2'	2:F:2:DT:H5'	2.00	0.90
1:G:25:CYS:HG	3:G:503:ZN:ZN	0.80	0.87
1:G:137:THR:HG22	1:G:139:LYS:H	1.40	0.86
1:C:154:GLU:HG2	1:C:286:VAL:HG22	1.56	0.85
1:E:21:PHE:HB3	1:E:25:CYS:O	1.76	0.85
1:A:242:VAL:HG11	1:A:261:GLY:CA	2.08	0.83
1:C:342:ARG:HG2	1:C:346:HIS:HB2	1.60	0.83
1:A:326:ASN:HD22	1:A:329:PHE:HB2	1.44	0.82
1:E:145:LYS:HE3	1:E:255:THR:O	1.80	0.82
1:A:239:VAL:CG1	1:A:242:VAL:HG12	2.10	0.81
1:E:266:ASN:HD22	1:E:308:ILE:HG23	1.44	0.81
1:G:237:THR:HB	1:G:259:VAL:HG12	1.63	0.80
1:G:173:LEU:HD21	1:G:201:ALA:HB2	1.62	0.80
1:E:372:ARG:HE	1:E:373:PRO:HD3	1.46	0.79
1:E:26:VAL:H	1:E:27:PRO:HD2	1.47	0.79
1:E:26:VAL:H	1:E:27:PRO:CD	1.96	0.78
1:E:293:VAL:HG11	1:G:293:VAL:HG11	1.64	0.78
1:E:239:VAL:HG12	1:E:242:VAL:HG12	1.64	0.77
1:A:358:PRO:HA	1:A:383:SER:HB3	1.66	0.76
1:A:242:VAL:HG11	1:A:261:GLY:HA2	1.65	0.75
1:G:267:GLN:O	1:G:278:ARG:NH2	2.20	0.75
1:A:71:ILE:CD1	1:A:147:LEU:HD21	2.15	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:154:GLU:HG2	1:C:286:VAL:CG2	2.17	0.75
1:C:159:PRO:O	1:C:162:SER:HB2	1.86	0.75
1:A:304:LEU:HD21	1:A:321:VAL:HG11	1.70	0.74
1:G:380:THR:HG23	1:G:381:ARG:HH21	1.51	0.74
2:B:1:DT:H1'	2:B:2:DT:O5'	1.88	0.74
1:G:72:LEU:O	1:G:76:LEU:HG	1.87	0.73
1:A:305:ILE:HD12	1:A:309:TYR:HD2	1.52	0.73
1:A:165:THR:HG22	1:A:195:LYS:HD2	1.69	0.73
1:G:357:PHE:O	1:G:383:SER:HB2	1.89	0.73
1:A:300:ALA:HB3	1:A:319:THR:O	1.89	0.73
1:C:165:THR:HG22	1:C:195:LYS:HD2	1.70	0.73
1:E:296:ARG:HD2	1:E:356:THR:HG21	1.71	0.72
1:E:242:VAL:HG11	1:E:261:GLY:CA	2.19	0.72
1:E:227:LEU:HB2	1:E:230:VAL:HG23	1.72	0.72
2:H:4:DT:C2'	2:H:5:DT:H72	2.10	0.71
1:G:84:LYS:HG3	1:G:135:GLN:HE22	1.54	0.71
1:G:331:LYS:HG3	1:G:332:GLY:H	1.55	0.71
1:A:278:ARG:NH1	5:A:693:HOH:O	2.23	0.70
1:G:358:PRO:HB3	1:G:384:GLN:HE21	1.56	0.70
1:A:181:LEU:CD1	1:A:238:PHE:HB2	2.21	0.69
1:A:247:PRO:HG3	1:A:276:VAL:HG11	1.74	0.69
1:E:335:ILE:HD12	1:E:349:ILE:HD13	1.75	0.68
1:E:335:ILE:HA	1:E:347:ARG:O	1.93	0.68
1:E:337:ALA:CB	1:E:369:SER:HB2	2.24	0.68
1:E:336:THR:HG21	1:E:341:ASP:HB2	1.75	0.68
1:G:372:ARG:O	1:G:376:VAL:HG23	1.94	0.67
1:E:212:VAL:HG23	1:E:213:LEU:HD12	1.75	0.67
1:C:300:ALA:HB1	1:C:321:VAL:HG12	1.75	0.67
1:E:352:ILE:HD13	1:E:381:ARG:HD2	1.77	0.67
1:C:364:LEU:HD21	1:C:370:LEU:HD11	1.75	0.67
1:C:233:SER:H	1:C:256:GLN:NE2	1.93	0.67
1:A:137:THR:O	1:A:252:ARG:NH2	2.28	0.67
1:G:168:LEU:HD21	1:G:238:PHE:HB3	1.77	0.67
1:G:160:PRO:O	1:G:312:TYR:OH	2.09	0.66
1:C:159:PRO:HA	1:C:264:ASP:OD2	1.94	0.66
1:G:145:LYS:NZ	1:G:255:THR:O	2.28	0.66
1:G:53:CYS:SG	1:G:56:CYS:HB3	2.36	0.65
1:A:233:SER:HB2	1:A:256:GLN:HB3	1.77	0.65
1:A:190:LEU:HG	1:A:194:ASN:HD21	1.59	0.65
1:E:293:VAL:CG1	1:G:293:VAL:HG11	2.26	0.65
1:E:108:LYS:HD3	1:E:109:VAL:H	1.62	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:3:VAL:HB	1:A:8:GLY:HA2	1.79	0.65
1:C:144:ASN:H	1:C:144:ASN:HD22	1.45	0.65
1:E:183:VAL:O	1:E:228:PRO:HD3	1.97	0.64
1:A:150:ALA:O	1:A:284:HIS:O	2.14	0.64
1:C:22:CYS:O	1:C:26:VAL:HG23	1.98	0.64
1:C:42:CYS:SG	1:C:44:HIS:HB2	2.37	0.64
1:A:334:VAL:HB	1:A:346:HIS:CD2	2.33	0.64
1:A:187:ALA:HB2	5:A:774:HOH:O	1.97	0.64
1:A:293:VAL:HG11	1:C:293:VAL:HG11	1.79	0.63
1:G:160:PRO:HD3	1:G:267:GLN:HE21	1.63	0.63
1:A:79:ILE:HG13	1:A:136:ASP:HB3	1.79	0.63
1:G:84:LYS:HG3	1:G:135:GLN:NE2	2.13	0.63
1:C:115:ASP:HB3	1:G:344:LEU:HD12	1.80	0.63
1:E:252:ARG:HG2	5:E:750:HOH:O	1.98	0.62
1:A:44:HIS:ND1	1:A:56:CYS:SG	2.72	0.62
1:G:382:ALA:HB2	1:G:386:LEU:HD13	1.80	0.62
1:E:238:PHE:HA	1:E:260:LYS:O	2.00	0.62
1:E:172:VAL:HG21	1:E:181:LEU:CD1	2.24	0.62
1:A:158:GLY:O	1:A:164:LYS:HE2	1.98	0.62
1:E:297:PHE:HE2	1:E:380:THR:O	1.81	0.62
1:C:182:VAL:HB	1:C:239:VAL:HG22	1.82	0.62
1:G:29:HIS:CD2	1:G:29:HIS:O	2.53	0.62
1:A:362:LEU:HD23	1:A:388:ILE:HG12	1.82	0.62
1:A:71:ILE:HD13	1:A:147:LEU:HD21	1.81	0.61
1:C:160:PRO:HG3	1:C:380:THR:HG21	1.81	0.61
1:A:355:CYS:O	1:A:381:ARG:NH1	2.30	0.61
1:A:339:HIS:CD2	2:B:3:DT:H5'	2.35	0.61
1:G:319:THR:HG23	1:G:384:GLN:HA	1.82	0.61
2:F:1:DT:H2'	2:F:2:DT:C5'	2.10	0.61
1:E:380:THR:HG22	1:E:381:ARG:NH2	2.16	0.61
1:C:160:PRO:HD3	1:C:267:GLN:NE2	2.16	0.60
1:G:159:PRO:O	1:G:162:SER:OG	2.11	0.60
1:E:158:GLY:O	1:E:164:LYS:HE2	2.02	0.60
1:G:359:VAL:HG13	1:G:385:GLU:HB3	1.83	0.60
1:G:25:CYS:SG	3:G:503:ZN:ZN	1.88	0.60
1:A:315:ALA:HB1	1:A:316:PRO:HD2	1.83	0.60
1:G:377:VAL:HA	1:G:380:THR:HG22	1.84	0.60
1:A:305:ILE:HD12	1:A:309:TYR:CD2	2.35	0.60
1:G:211:THR:HG21	1:G:215:PHE:HZ	1.67	0.60
1:E:251:ALA:HB2	1:E:281:TRP:CE2	2.35	0.60
1:A:21:PHE:CE2	1:A:29:HIS:HB2	2.36	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:160:PRO:HD3	1:G:267:GLN:NE2	2.17	0.59
1:A:323:PHE:O	1:A:325:PRO:HD3	2.01	0.59
1:E:156:VAL:HG13	1:E:289:GLU:O	2.02	0.59
1:G:376:VAL:O	1:G:380:THR:HB	2.01	0.59
1:E:321:VAL:O	1:E:321:VAL:HG13	2.01	0.59
1:G:183:VAL:HG22	1:G:226:ARG:O	2.02	0.59
1:C:183:VAL:HG21	1:C:189:MET:HG3	1.84	0.59
1:C:173:LEU:HD21	1:C:201:ALA:HB2	1.85	0.59
1:G:330:GLU:HG2	1:G:334:VAL:HG21	1.84	0.59
1:A:242:VAL:HG11	1:A:261:GLY:HA3	1.85	0.59
1:C:348:THR:HG22	1:C:350:ASP:H	1.68	0.59
1:C:394:GLN:O	1:C:398:LEU:HD12	2.02	0.59
1:C:59:SER:H	1:C:60:PRO:CD	2.16	0.59
1:A:316:PRO:HG3	1:C:295:HIS:CE1	2.38	0.58
1:C:124:CYS:SG	1:C:125:GLY:N	2.73	0.58
1:A:244:TYR:O	1:A:270:CYS:HB3	2.03	0.58
1:C:362:LEU:HD23	1:C:388:ILE:HG12	1.85	0.58
1:A:96:THR:CB	1:A:119:ASN:O	2.51	0.58
2:B:4:DT:H2"	2:B:5:DT:OP2	2.02	0.58
1:E:296:ARG:CD	1:E:356:THR:HG21	2.32	0.58
1:A:371:THR:HB	1:A:373:PRO:HD2	1.84	0.58
1:E:338:TYR:OH	1:E:369:SER:HA	2.03	0.58
1:C:2:ALA:HB3	1:C:20:TRP:CD1	2.39	0.58
1:C:364:LEU:CD2	1:C:370:LEU:HD11	2.33	0.57
1:A:395:LEU:O	1:A:399:LEU:HB2	2.04	0.57
1:E:230:VAL:HG22	2:F:7:DT:H5'	1.86	0.57
1:G:342:ARG:HG2	1:G:346:HIS:HB2	1.84	0.57
1:E:227:LEU:O	1:E:229:GLN:N	2.36	0.57
1:G:14:LYS:HB2	1:G:38:LEU:HD21	1.86	0.57
1:E:100:PRO:HA	1:E:112:VAL:HG22	1.86	0.57
1:A:264:ASP:OD2	1:A:267:GLN:NE2	2.33	0.57
1:E:3:VAL:O	1:E:20:TRP:HB2	2.04	0.57
1:G:380:THR:CG2	1:G:381:ARG:HH21	2.16	0.57
1:E:160:PRO:HB2	1:E:380:THR:CG2	2.35	0.57
1:E:149:ASN:ND2	1:E:257:GLY:HA2	2.21	0.56
1:A:326:ASN:HD21	1:A:329:PHE:HB2	1.70	0.56
1:E:249:ASP:O	1:E:252:ARG:HB3	2.05	0.56
1:C:227:LEU:HD13	2:D:7:DT:OP1	2.06	0.56
1:C:163:GLY:O	1:C:164:LYS:C	2.42	0.56
1:E:251:ALA:HB2	1:E:281:TRP:NE1	2.20	0.56
1:A:186:HIS:CE1	2:B:7:DT:OP1	2.58	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:240:ASP:OD1	1:C:241:GLU:HG2	2.05	0.56
1:G:240:ASP:O	1:G:241:GLU:C	2.43	0.56
1:E:293:VAL:HG11	1:G:293:VAL:CG1	2.33	0.56
1:C:144:ASN:HD22	1:C:144:ASN:N	2.02	0.55
1:C:359:VAL:HG22	1:C:385:GLU:HB2	1.87	0.55
1:C:184:PRO:HD2	1:C:188:SER:OG	2.07	0.55
2:F:6:DT:H2"	2:F:7:DT:H71	1.88	0.55
1:C:297:PHE:HB2	1:C:319:THR:OG1	2.07	0.55
1:E:156:VAL:HG22	1:E:288:LEU:HD12	1.89	0.55
1:C:59:SER:H	1:C:60:PRO:HD2	1.71	0.55
1:A:300:ALA:O	1:A:304:LEU:HD23	2.07	0.55
1:E:186:HIS:HA	1:E:189:MET:HB3	1.89	0.55
1:C:249:ASP:HA	1:C:252:ARG:HH11	1.72	0.55
1:C:4:CYS:HB3	1:C:9:ALA:H	1.73	0.54
1:A:158:GLY:O	1:A:164:LYS:CE	2.55	0.54
1:A:282:LEU:O	1:A:285:PHE:HB2	2.07	0.54
1:A:172:VAL:HG11	1:A:223:ILE:HG21	1.90	0.54
1:A:47:MET:O	1:A:62:GLN:HA	2.08	0.54
1:E:26:VAL:N	1:E:27:PRO:CD	2.66	0.54
1:C:160:PRO:O	1:C:312:TYR:OH	2.18	0.54
1:C:297:PHE:HE2	1:C:380:THR:HG1	1.54	0.54
1:C:352:ILE:HD12	1:C:381:ARG:HG3	1.89	0.53
1:E:147:LEU:O	1:E:150:ALA:HB3	2.08	0.53
1:E:158:GLY:HA3	1:E:164:LYS:HG2	1.91	0.53
1:A:395:LEU:HB3	1:A:401:PHE:HE2	1.73	0.53
1:C:227:LEU:HB3	1:C:228:PRO:HD2	1.91	0.53
1:G:137:THR:HG22	1:G:138:LEU:N	2.24	0.53
1:A:191:ASP:O	1:A:195:LYS:HG3	2.08	0.53
1:C:335:ILE:HG21	1:C:352:ILE:HD13	1.91	0.53
1:C:47:MET:O	1:C:63:MET:N	2.42	0.53
1:C:223:ILE:H	1:C:223:ILE:HD12	1.73	0.53
1:A:321:VAL:HG23	1:A:386:LEU:HB3	1.89	0.53
1:E:321:VAL:CG1	1:E:321:VAL:O	2.57	0.53
1:E:29:HIS:CG	1:E:29:HIS:O	2.62	0.53
1:G:331:LYS:HG3	1:G:332:GLY:N	2.24	0.53
1:A:244:TYR:HA	1:A:269:GLY:HA2	1.90	0.53
1:E:233:SER:H	1:E:256:GLN:NE2	2.08	0.52
1:E:316:PRO:HB2	1:G:166:PHE:HB3	1.90	0.52
1:G:182:VAL:HB	1:G:239:VAL:HG22	1.91	0.52
1:C:12:VAL:H	1:C:22:CYS:HA	1.74	0.52
1:G:4:CYS:HB3	1:G:9:ALA:H	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:50:SER:HB3	1:G:63:MET:HB3	1.91	0.52
1:E:358:PRO:HA	1:E:383:SER:CB	2.31	0.52
1:E:342:ARG:HD3	1:E:348:THR:HG23	1.91	0.52
1:C:283:ARG:HD2	1:C:284:HIS:NE2	2.23	0.52
1:G:4:CYS:HA	1:G:20:TRP:O	2.10	0.52
1:E:69:HIS:HB3	1:E:72:LEU:HB3	1.90	0.52
2:H:5:DT:H73	2:H:6:DT:H73	1.91	0.52
1:G:242:VAL:HG12	1:G:262:TYR:O	2.09	0.52
1:E:242:VAL:HG13	1:E:262:TYR:O	2.10	0.52
1:C:372:ARG:O	1:C:376:VAL:HG23	2.09	0.52
1:A:300:ALA:CB	1:A:319:THR:O	2.58	0.52
1:E:81:TYR:OH	2:F:3:DT:O4	2.22	0.52
1:E:160:PRO:HB2	1:E:380:THR:HG23	1.92	0.52
1:G:335:ILE:HG22	1:G:347:ARG:HB2	1.92	0.52
1:C:194:ASN:O	1:C:198:GLN:HG3	2.10	0.51
1:E:223:ILE:HG22	1:E:224:THR:N	2.25	0.51
1:A:145:LYS:HE2	1:A:255:THR:O	2.10	0.51
1:E:189:MET:HG2	1:E:215:PHE:HE2	1.75	0.51
1:A:233:SER:H	1:A:256:GLN:HE21	1.59	0.51
1:G:266:ASN:O	1:G:373:PRO:HA	2.10	0.51
1:G:341:ASP:OD2	1:G:366:THR:OG1	2.28	0.51
1:C:165:THR:CG2	1:C:195:LYS:HD2	2.39	0.51
1:E:208:PRO:HB2	1:E:211:THR:HB	1.92	0.51
1:A:137:THR:HG23	1:A:139:LYS:H	1.76	0.51
1:A:161:GLY:HA3	1:A:296:ARG:HB2	1.91	0.51
1:E:380:THR:CG2	1:E:381:ARG:NH2	2.73	0.51
1:A:98:SER:CB	1:A:99:PRO:HD3	2.41	0.51
1:E:296:ARG:NE	1:E:356:THR:HG21	2.26	0.51
1:A:190:LEU:HG	1:A:194:ASN:ND2	2.25	0.51
2:H:5:DT:C7	2:H:6:DT:H73	2.41	0.51
1:A:374:ARG:NH2	2:B:2:DT:O2	2.38	0.51
1:A:160:PRO:HD3	1:A:267:GLN:NE2	2.26	0.51
1:A:112:VAL:HA	1:A:120:ILE:O	2.11	0.51
1:A:82:GLY:O	1:A:83:SER:HB2	2.11	0.51
1:A:315:ALA:HB1	1:A:316:PRO:CD	2.41	0.50
1:C:376:VAL:O	1:C:380:THR:CG2	2.59	0.50
1:E:355:CYS:O	1:E:381:ARG:NH1	2.42	0.50
1:E:108:LYS:HD3	1:E:109:VAL:N	2.25	0.50
1:A:181:LEU:HD12	1:A:238:PHE:HB2	1.91	0.50
1:A:4:CYS:HB3	1:A:8:GLY:H	1.76	0.50
1:E:79:ILE:HD13	1:E:136:ASP:HB3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:304:LEU:HD12	1:A:379:VAL:HG11	1.93	0.50
1:A:360:VAL:HB	1:A:382:ALA:HB2	1.93	0.50
1:E:50:SER:OG	1:E:51:THR:N	2.45	0.50
1:G:168:LEU:O	1:G:172:VAL:HG23	2.11	0.50
1:C:98:SER:HB2	1:C:99:PRO:HD2	1.92	0.50
1:C:98:SER:HB2	1:C:99:PRO:CD	2.42	0.50
1:E:352:ILE:CD1	1:E:381:ARG:HD2	2.39	0.50
1:E:104:LYS:HG2	1:E:106:GLY:H	1.76	0.50
1:E:183:VAL:HG21	1:E:189:MET:HB2	1.94	0.49
1:E:251:ALA:HA	1:E:254:LEU:HD12	1.94	0.49
1:G:26:VAL:HG12	1:G:64:VAL:HG11	1.93	0.49
1:A:177:GLY:C	1:A:222:ASN:HD21	2.16	0.49
1:C:315:ALA:HB1	1:C:316:PRO:HD2	1.93	0.49
1:E:251:ALA:HB2	1:E:281:TRP:CZ2	2.46	0.49
1:E:149:ASN:HD22	1:E:257:GLY:HA2	1.77	0.49
1:C:38:LEU:HD23	5:C:620:HOH:O	2.11	0.49
1:C:377:VAL:O	1:C:381:ARG:HG2	2.13	0.49
1:C:376:VAL:O	1:C:380:THR:HG22	2.13	0.49
1:A:79:ILE:HG13	1:A:136:ASP:CB	2.42	0.49
1:E:83:SER:O	1:E:133:PRO:HG2	2.13	0.49
1:G:297:PHE:HE2	1:G:380:THR:O	1.96	0.49
1:A:299:ALA:N	1:A:314:PRO:HB3	2.28	0.49
1:A:183:VAL:O	1:A:228:PRO:HD3	2.13	0.48
2:F:2:DT:H2''	2:F:3:DT:O5'	2.14	0.48
1:G:377:VAL:O	1:G:381:ARG:HG2	2.13	0.48
1:A:169:VAL:HG22	1:A:196:LEU:HD13	1.95	0.48
1:E:372:ARG:N	1:E:373:PRO:HD2	2.29	0.48
1:A:294:CYS:SG	1:A:297:PHE:CE1	3.07	0.48
1:A:99:PRO:HB2	1:A:103:TYR:OH	2.13	0.48
1:C:6:VAL:O	1:C:144:ASN:ND2	2.46	0.48
1:C:226:ARG:HD3	1:C:232:THR:O	2.13	0.48
1:A:104:LYS:HG2	1:A:106:GLY:H	1.77	0.48
1:E:189:MET:HG2	1:E:215:PHE:CE2	2.49	0.48
1:A:226:ARG:NH2	1:A:230:VAL:O	2.37	0.48
1:G:358:PRO:HA	1:G:383:SER:HB3	1.94	0.48
1:G:62:GLN:NE2	1:G:64:VAL:O	2.47	0.48
1:A:134:LEU:HD12	1:A:134:LEU:O	2.12	0.48
1:C:122:PHE:C	1:C:124:CYS:H	2.17	0.48
1:E:316:PRO:HB3	1:G:167:HIS:CD2	2.49	0.48
1:E:110:VAL:HG12	1:E:123:GLY:HA3	1.95	0.48
1:G:245:PHE:HD2	1:G:249:ASP:HB3	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:167:HIS:HA	1:A:170:LYS:HE2	1.96	0.48
1:A:366:THR:O	1:A:369:SER:HB3	2.14	0.48
1:A:5:THR:O	1:A:143:VAL:HG23	2.13	0.48
1:E:247:PRO:HG3	1:E:276:VAL:HG11	1.95	0.48
1:E:181:LEU:HD23	1:E:192:CYS:HB3	1.95	0.47
1:C:168:LEU:CD1	1:C:240:ASP:HB2	2.44	0.47
1:A:72:LEU:O	1:A:76:LEU:HG	2.13	0.47
1:G:26:VAL:N	1:G:27:PRO:CD	2.77	0.47
1:G:282:LEU:HA	1:G:285:PHE:CD1	2.49	0.47
1:E:227:LEU:C	1:E:229:GLN:H	2.16	0.47
1:E:363:ARG:NH2	5:E:654:HOH:O	2.46	0.47
1:E:339:HIS:HA	1:E:342:ARG:HB2	1.95	0.47
1:G:374:ARG:O	1:G:378:ALA:CB	2.62	0.47
2:H:3:DT:H2''	2:H:4:DT:H5'	1.97	0.47
1:E:228:PRO:HB3	1:E:253:ILE:HD11	1.97	0.47
1:E:132:VAL:HG22	1:E:133:PRO:HD2	1.95	0.47
1:C:336:THR:HG22	1:C:363:ARG:HB2	1.96	0.47
1:A:71:ILE:CD1	1:A:147:LEU:CD2	2.88	0.47
1:C:321:VAL:HG13	1:C:321:VAL:O	2.14	0.47
1:A:352:ILE:HD13	1:A:381:ARG:HD2	1.97	0.47
1:A:172:VAL:HG12	1:A:223:ILE:HG12	1.97	0.47
1:G:350:ASP:OD1	1:G:374:ARG:NE	2.46	0.47
1:A:365:PRO:O	1:A:392:PHE:HE1	1.98	0.47
1:C:250:LEU:HA	1:C:253:ILE:HD12	1.97	0.47
1:A:380:THR:O	1:A:380:THR:HG23	2.15	0.47
1:E:335:ILE:HD13	1:E:381:ARG:HG3	1.96	0.47
1:G:363:ARG:NE	1:G:389:TYR:CD2	2.80	0.47
1:A:242:VAL:HG22	1:A:263:GLY:HA3	1.95	0.47
1:G:138:LEU:O	1:G:139:LYS:C	2.51	0.46
1:C:362:LEU:HD13	1:C:378:ALA:HB1	1.96	0.46
1:C:266:ASN:HD22	1:C:372:ARG:HH21	1.62	0.46
1:G:226:ARG:NH2	1:G:231:GLY:O	2.48	0.46
1:A:208:PRO:O	1:A:211:THR:HG22	2.16	0.46
1:C:265:LEU:HD21	1:C:282:LEU:HD12	1.96	0.46
1:G:32:HIS:HE2	1:G:80:ASP:CG	2.18	0.46
1:C:141:VAL:HG12	1:C:142:VAL:N	2.31	0.46
1:G:163:GLY:O	1:G:166:PHE:HB2	2.16	0.46
1:E:334:VAL:HG11	1:E:346:HIS:CE1	2.50	0.46
1:E:186:HIS:CE1	1:E:227:LEU:HD11	2.50	0.46
1:G:157:GLU:OE1	1:G:311:TYR:OH	2.26	0.46
1:C:371:THR:CG2	5:C:653:HOH:O	2.64	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:300:ALA:HB3	1:C:319:THR:O	2.16	0.46
1:E:171:ASP:HB3	1:E:238:PHE:CE2	2.51	0.46
2:F:1:DT:H1'	5:F:106:HOH:O	2.15	0.46
1:G:190:LEU:HG	5:G:684:HOH:O	2.15	0.46
1:C:191:ASP:O	1:C:195:LYS:HG3	2.15	0.45
1:E:337:ALA:HB3	1:E:369:SER:HB2	1.95	0.45
1:E:223:ILE:CG2	1:E:224:THR:N	2.79	0.45
1:A:202:ASP:O	1:A:221:GLY:HA3	2.16	0.45
1:A:166:PHE:O	1:A:170:LYS:HB2	2.17	0.45
1:G:217:ARG:HG3	5:G:658:HOH:O	2.17	0.45
1:A:250:LEU:O	1:A:251:ALA:C	2.55	0.45
1:E:69:HIS:HA	1:E:70:PRO:HD2	1.70	0.45
2:H:2:DT:H2''	2:H:3:DT:C6	2.52	0.45
1:A:182:VAL:HG13	1:A:226:ARG:HG3	1.97	0.45
1:C:185:THR:HB	2:D:6:DT:OP1	2.17	0.45
1:C:151:ALA:HA	1:C:284:HIS:O	2.16	0.45
1:G:363:ARG:HD2	1:G:389:TYR:HB3	1.98	0.45
1:A:351:SER:HB2	5:A:780:HOH:O	2.16	0.45
1:G:69:HIS:O	1:G:73:ASP:HB2	2.15	0.45
1:G:66:LYS:HB3	5:G:737:HOH:O	2.15	0.45
1:A:297:PHE:HE2	1:A:380:THR:O	2.00	0.45
1:C:99:PRO:HA	1:C:100:PRO:HD3	1.90	0.45
1:E:360:VAL:HB	1:E:382:ALA:HB2	1.99	0.45
1:A:35:THR:HB	1:A:46:ILE:HG22	1.98	0.45
1:A:372:ARG:O	1:A:376:VAL:HG23	2.16	0.44
1:E:363:ARG:HG2	1:E:389:TYR:HB3	1.99	0.44
1:A:248:VAL:O	1:A:252:ARG:HB2	2.17	0.44
1:E:163:GLY:O	1:E:164:LYS:C	2.54	0.44
1:G:215:PHE:CD1	1:G:215:PHE:N	2.84	0.44
1:A:99:PRO:HA	1:A:100:PRO:HD3	1.73	0.44
1:E:26:VAL:HG12	1:E:64:VAL:HG11	1.98	0.44
1:E:297:PHE:HD2	1:E:382:ALA:O	2.01	0.44
1:A:376:VAL:O	1:A:380:THR:HB	2.18	0.44
1:E:384:GLN:HA	1:E:384:GLN:HE21	1.82	0.44
1:G:297:PHE:CE2	1:G:380:THR:O	2.70	0.44
1:A:79:ILE:C	1:A:81:TYR:H	2.21	0.44
1:A:104:LYS:HG2	1:A:106:GLY:N	2.32	0.44
1:E:166:PHE:HB3	1:G:316:PRO:O	2.18	0.44
1:G:159:PRO:HB2	1:G:312:TYR:CZ	2.53	0.44
1:G:331:LYS:HG2	5:G:696:HOH:O	2.17	0.44
1:E:35:THR:O	1:E:45:ASP:HA	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:155:TYR:HB2	1:E:285:PHE:CD2	2.52	0.44
1:A:336:THR:HG21	1:A:341:ASP:HB2	2.00	0.44
1:C:286:VAL:O	1:C:286:VAL:CG2	2.65	0.43
1:G:183:VAL:HG23	1:G:227:LEU:HD23	1.98	0.43
1:A:100:PRO:HA	1:A:112:VAL:HG21	1.99	0.43
1:G:273:PRO:HG2	1:G:276:VAL:HG23	2.00	0.43
1:A:340:LYS:HB2	1:E:94:GLY:HA2	1.99	0.43
1:C:266:ASN:O	1:C:373:PRO:HA	2.18	0.43
1:C:5:THR:O	1:C:142:VAL:HA	2.18	0.43
1:E:266:ASN:ND2	1:E:308:ILE:HG23	2.23	0.43
1:A:71:ILE:HD13	1:A:147:LEU:CD2	2.47	0.43
1:C:233:SER:N	1:C:256:GLN:NE2	2.64	0.43
1:E:265:LEU:C	1:E:267:GLN:H	2.21	0.43
1:C:239:VAL:CG1	1:C:242:VAL:HG13	2.49	0.43
1:A:337:ALA:O	1:A:348:THR:HG23	2.18	0.43
1:C:379:VAL:HG21	1:C:399:LEU:HD21	2.01	0.43
1:E:228:PRO:HG3	1:E:245:PHE:HE2	1.84	0.43
1:E:160:PRO:HB2	1:E:380:THR:HG21	2.00	0.43
1:C:182:VAL:HG11	1:C:253:ILE:HG12	2.00	0.43
1:A:362:LEU:HD11	1:A:364:LEU:HD21	2.01	0.43
1:A:167:HIS:HE1	5:C:694:HOH:O	2.01	0.43
1:A:251:ALA:HB2	1:A:281:TRP:CZ2	2.53	0.43
1:G:308:ILE:HD12	1:G:372:ARG:HH21	1.83	0.43
1:A:13:ALA:HB1	1:A:35:THR:HG22	1.99	0.43
1:G:168:LEU:CD2	1:G:238:PHE:HB3	2.46	0.43
1:E:108:LYS:CD	1:E:109:VAL:H	2.29	0.43
1:C:17:CYS:SG	1:C:138:LEU:HD12	2.58	0.43
1:E:349:ILE:HA	1:E:352:ILE:HD11	1.99	0.43
1:A:186:HIS:CE1	1:A:208:PRO:HB3	2.54	0.43
1:A:172:VAL:HG11	1:A:223:ILE:CG2	2.48	0.43
1:E:110:VAL:HG12	1:E:124:CYS:H	1.83	0.43
1:C:202:ASP:O	1:C:221:GLY:HA3	2.18	0.43
1:G:372:ARG:HD3	5:G:638:HOH:O	2.19	0.43
1:A:46:ILE:C	1:A:48:TYR:H	2.22	0.43
2:H:7:DT:O4'	2:H:7:DT:O2	2.36	0.43
1:G:250:LEU:HA	1:G:253:ILE:HD12	2.00	0.43
1:G:321:VAL:HB	1:G:386:LEU:HB3	2.00	0.42
2:F:4:DT:O3'	2:F:5:DT:O2	2.37	0.42
1:G:5:THR:O	1:G:142:VAL:HA	2.19	0.42
1:C:244:TYR:HA	1:C:269:GLY:HA2	2.01	0.42
1:E:166:PHE:HB3	1:G:316:PRO:HB2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:60:PRO:HD3	5:G:757:HOH:O	2.19	0.42
1:A:17:CYS:SG	1:A:29:HIS:HE1	2.42	0.42
1:G:330:GLU:HG2	1:G:334:VAL:CG2	2.49	0.42
1:C:0:MET:O	1:C:1:SER:OG	2.20	0.42
1:C:164:LYS:HG2	1:C:165:THR:N	2.34	0.42
1:E:160:PRO:CB	1:E:380:THR:HG21	2.50	0.42
1:G:185:THR:HG1	1:G:188:SER:H	1.67	0.42
1:G:11:PRO:HA	1:G:22:CYS:HA	2.01	0.42
1:C:371:THR:HG21	5:C:653:HOH:O	2.19	0.42
1:C:70:PRO:HB2	1:C:71:ILE:HD12	2.01	0.42
1:E:371:THR:HB	1:E:373:PRO:HD2	2.01	0.42
1:G:47:MET:HA	1:G:64:VAL:HB	2.02	0.42
1:E:201:ALA:O	1:E:203:PRO:HD3	2.19	0.42
1:C:137:THR:HG23	1:C:139:LYS:H	1.85	0.42
1:E:47:MET:HA	1:E:64:VAL:HB	2.01	0.42
1:G:26:VAL:CG1	1:G:64:VAL:HG11	2.50	0.42
1:G:190:LEU:HD21	1:G:214:ASP:O	2.20	0.42
1:C:344:LEU:O	1:C:346:HIS:HD2	2.03	0.42
1:C:363:ARG:HD2	1:C:389:TYR:HD2	1.84	0.42
1:G:32:HIS:NE2	1:G:80:ASP:OD2	2.52	0.42
1:E:15:SER:HB3	1:E:35:THR:CG2	2.50	0.42
1:E:155:TYR:HB2	1:E:285:PHE:HD2	1.84	0.42
1:A:239:VAL:HG12	1:A:242:VAL:CG1	2.31	0.42
1:A:67:VAL:HA	1:A:68:PRO:HD3	1.92	0.42
1:C:366:THR:HA	1:C:367:PRO:HD3	1.89	0.42
1:C:188:SER:HA	1:C:191:ASP:HB2	2.02	0.42
1:C:352:ILE:CD1	1:C:381:ARG:HG3	2.50	0.42
1:C:317:HIS:HB2	1:C:318:THR:H	1.53	0.42
1:A:21:PHE:CZ	1:A:29:HIS:ND1	2.88	0.41
1:C:351:SER:OG	2:D:4:DT:OP1	2.36	0.41
2:H:4:DT:H2"	2:H:5:DT:C5	2.54	0.41
1:A:323:PHE:HZ	1:A:401:PHE:CZ	2.38	0.41
1:E:363:ARG:HD2	1:E:389:TYR:CD2	2.55	0.41
1:G:357:PHE:C	1:G:383:SER:HB2	2.40	0.41
1:A:350:ASP:OD2	2:B:3:DT:H1'	2.20	0.41
1:G:282:LEU:O	1:G:285:PHE:HB2	2.20	0.41
1:C:389:TYR:CE1	1:C:391:PRO:HD3	2.55	0.41
1:A:27:PRO:HB2	1:A:67:VAL:HG22	2.02	0.41
1:C:132:VAL:HG22	1:C:133:PRO:HD2	2.02	0.41
1:G:84:LYS:HG2	1:G:132:VAL:HG23	2.02	0.41
1:G:47:MET:HG2	1:G:48:TYR:N	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:171:ASP:OD2	1:C:262:TYR:OH	2.33	0.41
1:E:98:SER:H	1:E:99:PRO:CD	2.33	0.41
1:A:152:ALA:HB1	1:A:258:ARG:HD3	2.02	0.41
1:C:286:VAL:HG23	1:C:286:VAL:O	2.20	0.41
1:G:278:ARG:HD2	5:G:706:HOH:O	2.19	0.41
1:C:372:ARG:N	1:C:373:PRO:CD	2.84	0.41
1:A:297:PHE:O	1:A:315:ALA:HB3	2.20	0.41
1:E:49:ARG:HG2	1:E:50:SER:N	2.35	0.41
1:C:395:LEU:O	1:C:399:LEU:HB2	2.20	0.41
1:G:320:LYS:HD2	1:G:322:VAL:HG23	2.03	0.41
1:G:148:LYS:NZ	1:G:149:ASN:OD1	2.53	0.41
1:A:147:LEU:HD22	1:A:281:TRP:CZ3	2.55	0.41
1:E:171:ASP:O	1:E:175:VAL:HG23	2.21	0.41
1:A:298:GLY:HA2	1:A:315:ALA:H	1.85	0.41
1:A:380:THR:O	1:A:380:THR:CG2	2.67	0.41
1:C:331:LYS:HD2	1:C:359:VAL:HG11	2.02	0.41
1:E:49:ARG:CG	1:E:50:SER:N	2.84	0.41
1:C:363:ARG:HD2	1:C:389:TYR:CD2	2.56	0.41
1:E:264:ASP:OD2	1:E:267:GLN:HG3	2.21	0.41
1:C:69:HIS:HA	1:C:70:PRO:HD3	1.93	0.41
1:G:172:VAL:HG22	1:G:238:PHE:CD1	2.56	0.41
1:A:335:ILE:HD12	1:A:349:ILE:HG12	2.03	0.41
1:C:310:PRO:O	1:C:312:TYR:N	2.54	0.40
1:C:228:PRO:HG3	1:C:245:PHE:HE2	1.86	0.40
1:G:155:TYR:HD2	1:G:287:SER:OG	2.03	0.40
1:E:47:MET:HG3	1:E:62:GLN:HG3	2.03	0.40
1:C:268:LEU:HD23	1:C:373:PRO:HB2	2.03	0.40
1:G:246:SER:HA	1:G:270:CYS:SG	2.61	0.40
1:A:132:VAL:HG22	1:A:133:PRO:HD2	2.04	0.40
1:C:10:ALA:HA	1:C:11:PRO:HD3	1.84	0.40
1:C:273:PRO:HG2	1:C:276:VAL:HG23	2.03	0.40
1:C:109:VAL:O	1:C:109:VAL:HG23	2.22	0.40
1:C:264:ASP:OD1	1:C:266:ASN:HB2	2.21	0.40
1:C:240:ASP:OD1	1:C:241:GLU:N	2.54	0.40
1:C:141:VAL:HG13	1:C:251:ALA:HB1	2.04	0.40
1:C:370:LEU:HD23	1:C:374:ARG:HD2	2.03	0.40
1:C:283:ARG:HD2	1:C:284:HIS:CE1	2.57	0.40
1:A:335:ILE:HA	1:A:347:ARG:O	2.22	0.40
1:C:180:THR:O	1:C:237:THR:HA	2.20	0.40
1:G:271:VAL:HG22	5:G:730:HOH:O	2.21	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	383/423 (90%)	316 (82%)	54 (14%)	13 (3%)	5	9
1	C	390/423 (92%)	324 (83%)	49 (13%)	17 (4%)	3	5
1	E	386/423 (91%)	332 (86%)	45 (12%)	9 (2%)	8	17
1	G	385/423 (91%)	326 (85%)	48 (12%)	11 (3%)	6	12
All	All	1544/1692 (91%)	1298 (84%)	196 (13%)	50 (3%)	5	10

All (50) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	54	THR
1	A	98	SER
1	A	159	PRO
1	A	165	THR
1	C	118	GLY
1	C	143	VAL
1	G	393	ASP
1	C	1	SER
1	C	59	SER
1	C	106	GLY
1	C	141	VAL
1	C	311	TYR
1	C	343	GLY
1	E	266	ASN
1	G	120	ILE
1	G	123	GLY
1	G	221	GLY
1	A	46	ILE
1	A	299	ALA
1	C	133	PRO
1	C	314	PRO
1	E	50	SER

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Mol	Chain	Res	Type
1	E	287	SER
1	G	53	CYS
1	G	109	VAL
1	G	184	PRO
1	G	350	ASP
1	A	17	CYS
1	A	30	ALA
1	A	314	PRO
1	C	184	PRO
1	G	203	PRO
1	C	138	LEU
1	E	26	VAL
1	G	80	ASP
1	A	251	ALA
1	C	58	GLY
1	E	70	PRO
1	E	100	PRO
1	E	279	ASN
1	G	57	GLU
1	C	113	VAL
1	C	310	PRO
1	E	98	SER
1	E	228	PRO
1	C	358	PRO
1	A	316	PRO
1	A	125	GLY
1	C	142	VAL
1	A	158	GLY

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	300/352 (85%)	264 (88%)	36 (12%)	6	13
1	C	301/352 (86%)	265 (88%)	36 (12%)	6	13
1	E	298/352 (85%)	265 (89%)	33 (11%)	8	15

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	300/352 (85%)	273 (91%)	27 (9%)	12	24
All	All	1199/1408 (85%)	1067 (89%)	132 (11%)	8	16

All (132) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	CYS
1	A	35	THR
1	A	51	THR
1	A	53	CYS
1	A	56	CYS
1	A	64	VAL
1	A	89	LEU
1	A	129	HIS
1	A	134	LEU
1	A	136	ASP
1	A	137	THR
1	A	144	ASN
1	A	147	LEU
1	A	153	SER
1	A	168	LEU
1	A	182	VAL
1	A	196	LEU
1	A	214	ASP
1	A	226	ARG
1	A	236	GLU
1	A	246	SER
1	A	270	CYS
1	A	283	ARG
1	A	296	ARG
1	A	297	PHE
1	A	308	ILE
1	A	326	ASN
1	A	331	LYS
1	A	336	THR
1	A	355	CYS
1	A	361	THR
1	A	371	THR
1	A	380	THR
1	A	384	GLN
1	A	386	LEU
1	A	398	LEU

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Mol	Chain	Res	Type
1	C	12	VAL
1	C	22	CYS
1	C	42	CYS
1	C	44	HIS
1	C	51	THR
1	C	55	MET
1	C	64	VAL
1	C	75	LEU
1	C	88	THR
1	C	112	VAL
1	C	122	PHE
1	C	132	VAL
1	C	137	THR
1	C	144	ASN
1	C	162	SER
1	C	164	LYS
1	C	188	SER
1	C	220	SER
1	C	234	GLU
1	C	242	VAL
1	C	256	GLN
1	C	286	VAL
1	C	293	VAL
1	C	297	PHE
1	C	304	LEU
1	C	317	HIS
1	C	318	THR
1	C	350	ASP
1	C	356	THR
1	C	358	PRO
1	C	359	VAL
1	C	362	LEU
1	C	371	THR
1	C	380	THR
1	C	393	ASP
1	C	395	LEU
1	E	4	CYS
1	E	12	VAL
1	E	35	THR
1	E	107	HIS
1	E	112	VAL
1	E	132	VAL

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Mol	Chain	Res	Type
1	E	145	LYS
1	E	147	LEU
1	E	165	THR
1	E	166	PHE
1	E	176	VAL
1	E	181	LEU
1	E	210	TYR
1	E	220	SER
1	E	224	THR
1	E	230	VAL
1	E	233	SER
1	E	242	VAL
1	E	256	GLN
1	E	258	ARG
1	E	270	CYS
1	E	297	PHE
1	E	306	LYS
1	E	312	TYR
1	E	336	THR
1	E	339	HIS
1	E	341	ASP
1	E	348	THR
1	E	355	CYS
1	E	357	PHE
1	E	363	ARG
1	E	384	GLN
1	E	398	LEU
1	G	36	THR
1	G	51	THR
1	G	66	LYS
1	G	75	LEU
1	G	88	THR
1	G	132	VAL
1	G	145	LYS
1	G	156	VAL
1	G	164	LYS
1	G	168	LEU
1	G	192	CYS
1	G	215	PHE
1	G	226	ARG
1	G	287	SER
1	G	291	LEU

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Mol	Chain	Res	Type
1	G	304	LEU
1	G	306	LYS
1	G	312	TYR
1	G	321	VAL
1	G	322	VAL
1	G	340	LYS
1	G	371	THR
1	G	380	THR
1	G	383	SER
1	G	385	GLU
1	G	395	LEU
1	G	399	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	HIS
1	A	74	HIS
1	A	144	ASN
1	A	186	HIS
1	A	194	ASN
1	A	222	ASN
1	A	256	GLN
1	A	266	ASN
1	A	326	ASN
1	A	384	GLN
1	C	74	HIS
1	C	144	ASN
1	C	256	GLN
1	C	266	ASN
1	C	267	GLN
1	C	346	HIS
1	E	149	ASN
1	E	186	HIS
1	E	222	ASN
1	E	256	GLN
1	E	266	ASN
1	E	384	GLN
1	G	62	GLN
1	G	69	HIS
1	G	135	GLN
1	G	144	ASN

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Mol	Chain	Res	Type
1	G	167	HIS
1	G	186	HIS
1	G	198	GLN
1	G	256	GLN
1	G	267	GLN
1	G	384	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 19 ligands modelled in this entry, 19 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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	395/423 (93%)	0.80	38 (9%) 10 8	26, 48, 69, 83	0
1	C	398/423 (94%)	0.79	37 (9%) 11 8	27, 49, 67, 87	1 (0%)
1	E	396/423 (93%)	0.89	42 (10%) 8 6	30, 50, 72, 85	0
1	G	393/423 (92%)	0.85	45 (11%) 6 5	29, 50, 73, 94	0
2	B	7/7 (100%)	1.09	1 (14%) 4 2	54, 57, 67, 73	3 (42%)
2	D	7/7 (100%)	1.41	2 (28%) 1 0	60, 65, 67, 67	4 (57%)
2	F	7/7 (100%)	0.88	1 (14%) 4 2	54, 60, 63, 71	4 (57%)
2	H	7/7 (100%)	0.82	1 (14%) 4 2	49, 54, 62, 64	4 (57%)
All	All	1610/1720 (93%)	0.84	167 (10%) 8 6	26, 50, 70, 94	16 (0%)

All (167) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	335	ILE	6.8
1	E	62	GLN	5.5
1	C	297	PHE	5.0
1	A	177	GLY	4.9
1	C	344	LEU	4.9
1	G	55	MET	4.8
1	E	48	TYR	4.7
1	G	307	GLY	4.7
1	E	215	PHE	4.6
1	A	109	VAL	4.6
1	E	116	VAL	4.5
1	E	112	VAL	4.4
1	G	388	ILE	4.4
1	A	91	VAL	4.3
1	C	337	ALA	4.1
1	G	56	CYS	4.1

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Mol	Chain	Res	Type	RSRZ
1	C	343	GLY	4.1
1	A	82	GLY	4.0
1	C	60	PRO	4.0
1	C	348	THR	4.0
1	E	179	ALA	3.9
1	C	128	SER	3.9
1	G	51	THR	3.9
1	E	225	VAL	3.9
1	G	132	VAL	3.8
1	E	307	GLY	3.8
1	G	112	VAL	3.8
1	A	307	GLY	3.8
2	D	1	DT	3.7
1	G	61	LYS	3.7
1	C	109	VAL	3.5
1	C	336	THR	3.5
1	C	108	LYS	3.5
1	E	55	MET	3.4
1	E	93	ASP	3.3
1	E	105	VAL	3.3
1	G	113	VAL	3.3
1	E	124	CYS	3.3
1	G	108	LYS	3.3
1	A	188	SER	3.3
1	G	15	SER	3.2
1	A	111	ALA	3.2
1	G	31	GLY	3.2
1	C	132	VAL	3.2
1	E	321	VAL	3.2
1	A	223	ILE	3.2
2	H	1	DT	3.2
1	E	371	THR	3.2
1	E	129	HIS	3.1
1	E	108	LYS	3.1
1	G	19	GLY	3.1
1	G	87	LEU	3.0
1	G	50	SER	3.0
1	G	192	CYS	3.0
1	G	65	PRO	3.0
1	A	225	VAL	3.0
1	C	9	ALA	3.0
1	A	220	SER	3.0

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Mol	Chain	Res	Type	RSRZ
1	C	105	VAL	3.0
1	G	53	CYS	3.0
1	C	10	ALA	2.9
1	G	294	CYS	2.9
1	E	365	PRO	2.9
1	E	3	VAL	2.9
1	A	389	TYR	2.9
1	G	205	PHE	2.9
1	E	325	PRO	2.9
1	C	287	SER	2.9
1	G	59	SER	2.9
1	C	59	SER	2.8
1	G	152	ALA	2.8
1	A	122	PHE	2.8
1	E	109	VAL	2.8
1	E	213	LEU	2.8
1	E	160	PRO	2.8
1	E	82	GLY	2.8
1	G	29	HIS	2.8
1	G	195	LYS	2.7
1	E	71	ILE	2.7
1	C	217	ARG	2.7
1	A	237	THR	2.7
1	A	16	ALA	2.7
1	A	92	ALA	2.7
1	E	324	VAL	2.6
1	C	50	SER	2.6
1	A	131	ALA	2.6
1	E	188	SER	2.6
1	E	387	TYR	2.6
1	C	12	VAL	2.6
1	G	365	PRO	2.6
1	G	84	LYS	2.6
1	C	70	PRO	2.6
1	G	124	CYS	2.5
1	A	110	VAL	2.5
1	E	175	VAL	2.5
1	A	103	TYR	2.5
1	C	315	ALA	2.5
2	F	1	DT	2.5
1	C	308	ILE	2.5
1	G	217	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
1	E	216	PRO	2.5
1	A	105	VAL	2.4
1	C	51	THR	2.4
1	E	275	SER	2.4
1	A	398	LEU	2.4
1	E	308	ILE	2.4
1	G	345	GLY	2.4
1	G	207	VAL	2.4
1	E	369	SER	2.4
1	A	6	VAL	2.3
1	A	54	THR	2.3
1	A	98	SER	2.3
1	A	96	THR	2.3
1	C	65	PRO	2.3
1	G	126	PRO	2.3
1	C	398	LEU	2.3
1	E	240	ASP	2.3
1	A	209	LYS	2.3
1	C	386	LEU	2.3
1	A	23	GLY	2.2
2	D	6	DT	2.2
1	C	113	VAL	2.2
1	E	182	VAL	2.2
1	G	11	PRO	2.2
1	E	58	GLY	2.2
1	G	117	GLY	2.2
1	C	212	VAL	2.2
1	G	203	PRO	2.2
1	A	44	HIS	2.2
1	A	146	ALA	2.2
1	E	311	TYR	2.2
1	G	83	SER	2.2
1	A	106	GLY	2.2
1	C	123	GLY	2.2
1	C	106	GLY	2.1
2	B	7	DT	2.1
1	G	46	ILE	2.1
1	A	93	ASP	2.1
1	G	326	ASN	2.1
1	E	100	PRO	2.1
1	A	338	TYR	2.1
1	A	299	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	E	268	LEU	2.1
1	C	39	PHE	2.1
1	E	290	PRO	2.1
1	G	159	PRO	2.1
1	G	224	THR	2.1
1	G	28	TYR	2.1
1	A	57	GLU	2.1
1	C	200	GLY	2.1
1	E	111	ALA	2.1
1	G	173	LEU	2.1
1	E	329	PHE	2.1
1	G	111	ALA	2.1
1	E	54	THR	2.0
1	C	19	GLY	2.0
1	A	51	THR	2.0
1	G	116	VAL	2.0
1	A	401	PHE	2.0
1	A	43	GLY	2.0
1	C	201	ALA	2.0
1	G	30	ALA	2.0
1	G	386	LEU	2.0
1	C	172	VAL	2.0
1	A	214	ASP	2.0
1	C	107	HIS	2.0
1	A	322	VAL	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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
4	CA	C	505	1/1	0.91	0.20	-0.00	58,58,58,58	0
3	ZN	E	503	1/1	0.98	0.18	-0.00	62,62,62,62	0
4	CA	C	504	1/1	0.99	0.21	-0.22	46,46,46,46	0
3	ZN	A	501	1/1	0.98	0.18	-0.33	53,53,53,53	0
4	CA	E	505	1/1	0.98	0.22	-0.60	61,61,61,61	0
4	CA	G	504	1/1	0.95	0.15	-0.99	38,38,38,38	0
3	ZN	G	502	1/1	0.95	0.13	-1.13	78,78,78,78	0
3	ZN	E	502	1/1	0.95	0.09	-1.59	69,69,69,69	0
3	ZN	C	501	1/1	0.99	0.14	-1.61	42,42,42,42	0
3	ZN	A	503	1/1	0.99	0.12	-1.64	46,46,46,46	0
3	ZN	G	501	1/1	0.94	0.11	-1.66	52,52,52,52	0
3	ZN	G	503	1/1	0.95	0.12	-2.13	64,64,64,64	0
3	ZN	E	501	1/1	0.99	0.11	-2.46	57,57,57,57	0
3	ZN	C	503	1/1	0.99	0.11	-3.49	41,41,41,41	0
4	CA	A	504	1/1	0.83	0.21	-	58,58,58,58	0
3	ZN	A	502	1/1	0.93	0.11	-	73,73,73,73	0
4	CA	A	505	1/1	0.88	0.16	-	77,77,77,77	0
4	CA	E	504	1/1	0.76	0.14	-	65,65,65,65	0
3	ZN	C	502	1/1	0.97	0.10	-	83,83,83,83	0

6.5 Other polymers

There are no such residues in this entry.