



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 27, 2022 – 10:05 AM EST

PDB ID : 7MLB
Title : Crystal structure of Thermus thermophilus transcription initiation complex with 5nt RNA
Authors : Liu, Y.; Ebright, R.H.
Deposited on : 2021-04-28
Resolution : 3.60 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

<https://www.wwpdb.org/validation/2017/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
Xtriage (Phenix) : 1.13
EDS : 2.26
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.26

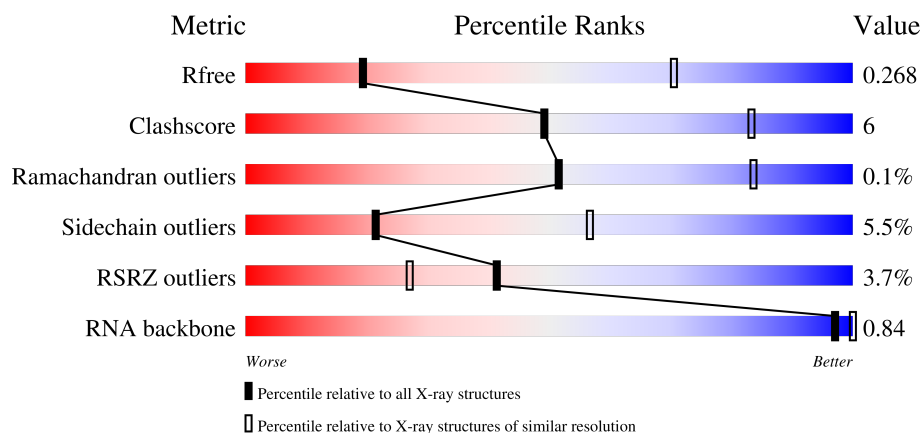
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.60 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}	130704	1257 (3.70-3.50)
Clashscore	141614	1353 (3.70-3.50)
Ramachandran outliers	138981	1307 (3.70-3.50)
Sidechain outliers	138945	1307 (3.70-3.50)
RSRZ outliers	127900	1161 (3.70-3.50)
RNA backbone	3102	1017 (4.20-3.00)

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	315	
1	B	315	
2	C	1119	
3	D	1524	

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Mol	Chain	Length	Quality of chain
4	E	99	
5	F	443	
6	G	21	
7	H	27	
8	I	5	

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
9	MG	B	2001	-	-	-	X

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 28400 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 DNA-directed RNA polymerase subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	226	Total	C	N	O	S	0	0	0
			1782	1138	310	332	2			
1	B	222	Total	C	N	O	S	0	0	0
			1750	1118	304	326	2			

- Molecule 2 is a protein called DNA-directed RNA polymerase subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	1111	Total	C	N	O	S	0	0	0
			8770	5548	1564	1634	24			

- Molecule 3 is a protein called DNA-directed RNA polymerase subunit beta'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	1486	Total	C	N	O	S	0	0	0
			11738	7441	2067	2195	35			

- Molecule 4 is a protein called DNA-directed RNA polymerase subunit omega.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	E	94	Total	C	N	O	S	0	0	0
			761	486	132	139	4			

- Molecule 5 is a protein called RNA polymerase sigma factor SigA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	F	346	Total	C	N	O	S	0	0	0
			2807	1770	509	524	4			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	-19	MET	-	initiating methionine	UNP Q5SKW1
F	-18	GLY	-	expression tag	UNP Q5SKW1
F	-17	SER	-	expression tag	UNP Q5SKW1
F	-16	SER	-	expression tag	UNP Q5SKW1
F	-15	HIS	-	expression tag	UNP Q5SKW1
F	-14	HIS	-	expression tag	UNP Q5SKW1
F	-13	HIS	-	expression tag	UNP Q5SKW1
F	-12	HIS	-	expression tag	UNP Q5SKW1
F	-11	HIS	-	expression tag	UNP Q5SKW1
F	-10	HIS	-	expression tag	UNP Q5SKW1
F	-9	SER	-	expression tag	UNP Q5SKW1
F	-8	SER	-	expression tag	UNP Q5SKW1
F	-7	GLY	-	expression tag	UNP Q5SKW1
F	-6	LEU	-	expression tag	UNP Q5SKW1
F	-5	VAL	-	expression tag	UNP Q5SKW1
F	-4	PRO	-	expression tag	UNP Q5SKW1
F	-3	ARG	-	expression tag	UNP Q5SKW1
F	-2	GLY	-	expression tag	UNP Q5SKW1
F	-1	SER	-	expression tag	UNP Q5SKW1
F	0	HIS	-	expression tag	UNP Q5SKW1

- Molecule 6 is a DNA chain called DNA (5'-D(P*CP*TP*GP*CP*AP*TP*CP*CP*CP*TP*GP*AP*GP*TP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	G	15	Total	C	N	O	P	0	0	0
			305	145	53	92	15			

- Molecule 7 is a DNA chain called DNA (5'-D(*TP*AP*TP*AP*AP*TP*GP*GP*AP*GP*GP*GP*AP*TP*GP*CP*AP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	H	18	Total	C	N	O	P	0	0	0
			376	179	76	104	17			

- Molecule 8 is a RNA chain called RNA (5'-R(P*AP*CP*UP*CP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	I	5	Total	C	N	O	P	0	0	0
			104	47	18	34	5			

- Molecule 9 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	B	1	Total 1	Mg 1	0	0
9	D	3	Total 3	Mg 3	0	0
9	F	1	Total 1	Mg 1	0	0

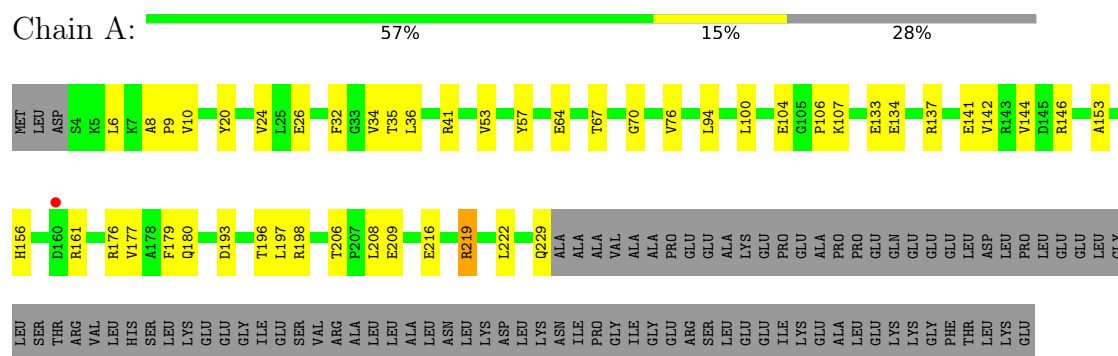
- Molecule 10 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	D	2	Total 2	Zn 2	0	0

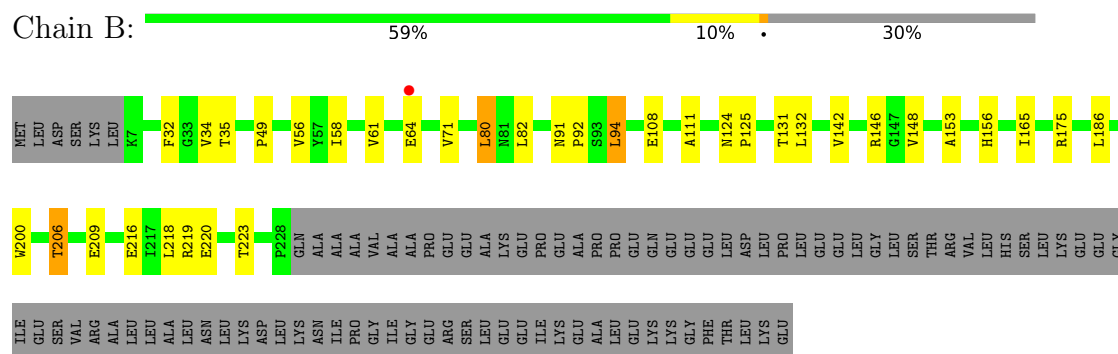
3 Residue-property plots [i](#)

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.

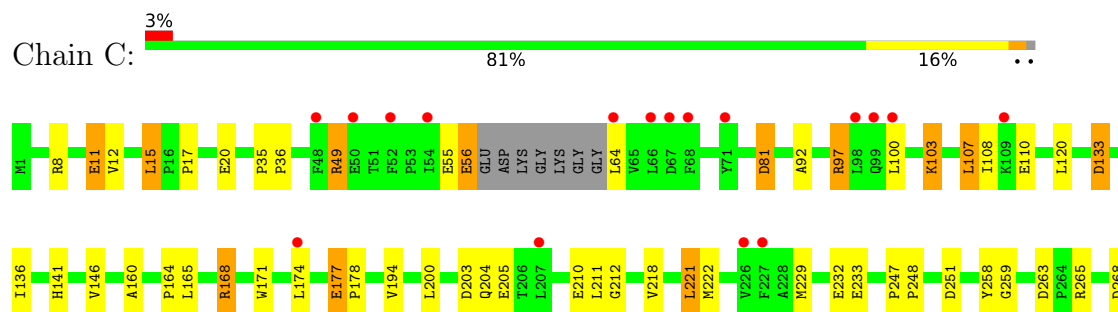
- Molecule 1: DNA-directed RNA polymerase subunit alpha

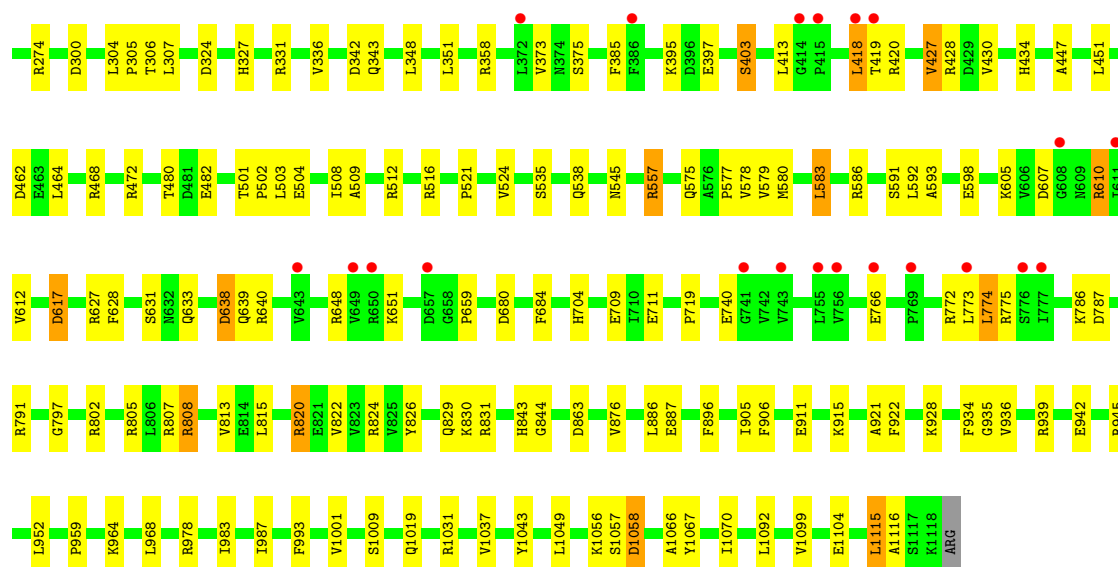


- Molecule 1: DNA-directed RNA polymerase subunit alpha

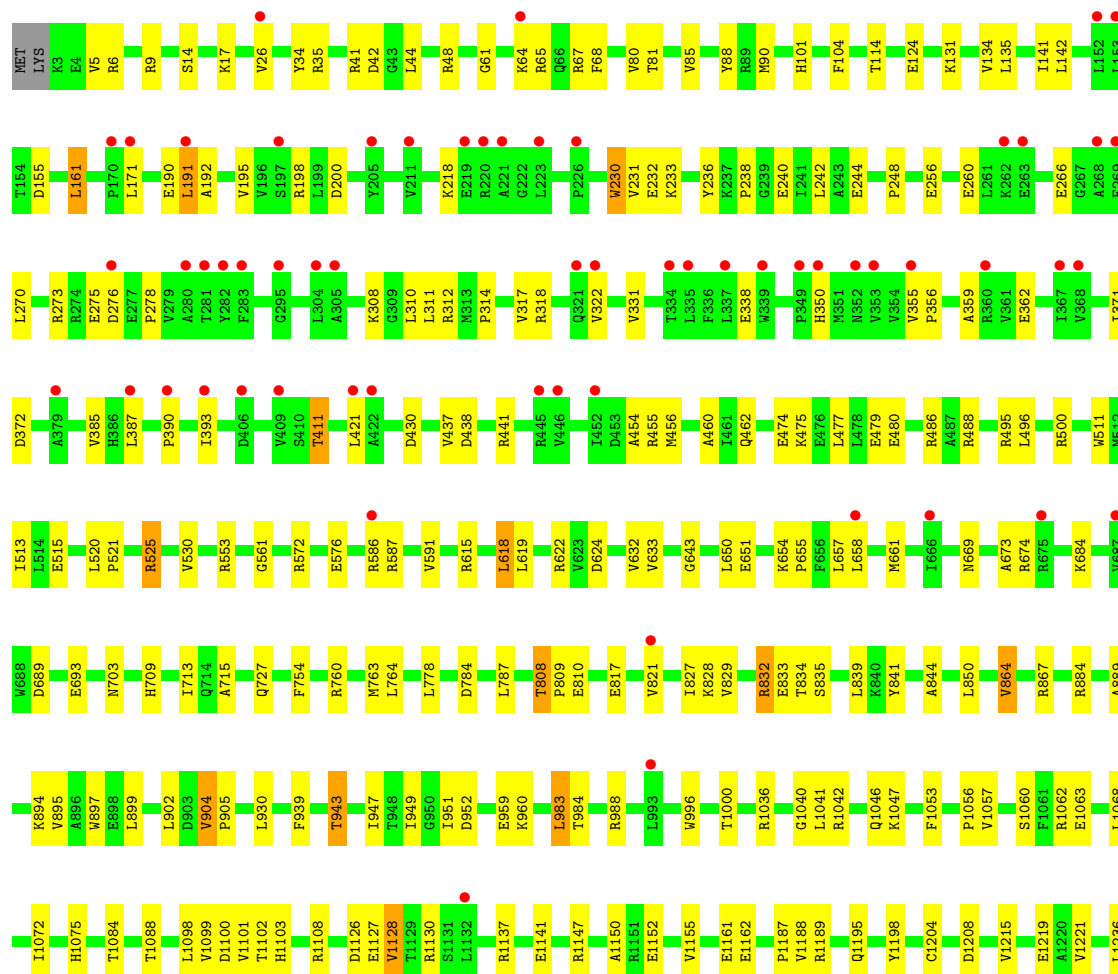
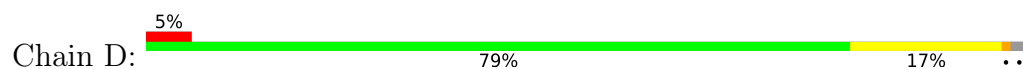


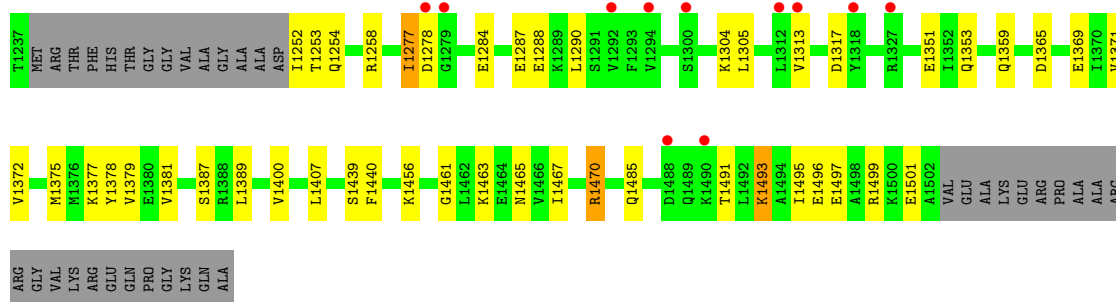
- Molecule 2: DNA-directed RNA polymerase subunit beta



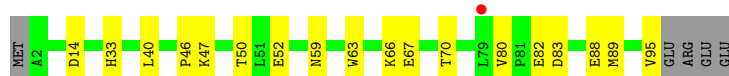
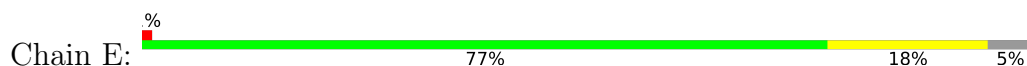


• Molecule 3: DNA-directed RNA polymerase subunit beta'

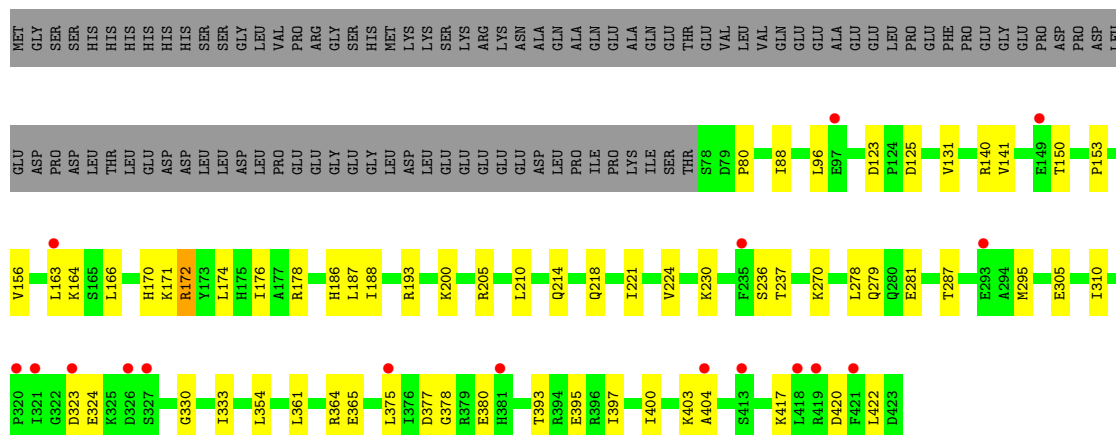




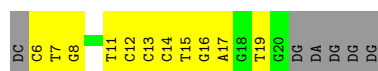
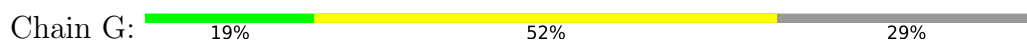
- Molecule 4: DNA-directed RNA polymerase subunit omega



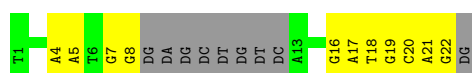
- Molecule 5: RNA polymerase sigma factor SigA



- Molecule 6: DNA (5'-D(P*CP*TP*GP*CP*AP*TP*CP*CP*CP*TP*GP*AP*GP*TP*G)-3')



- Molecule 7: DNA (5'-D(*TP*AP*TP*AP*AP*TP*GP*GP*AP*GP*GP*GP*AP*TP*GP*CP*AP*G)-3')



- Molecule 8: RNA (5'-R(P*AP*CP*UP*CP*A)-3')

Chain I:  60% 40%



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	186.17Å 104.27Å 299.02Å 90.00° 97.95° 90.00°	Depositor
Resolution (Å)	44.62 – 3.60 49.36 – 3.60	Depositor EDS
% Data completeness (in resolution range)	98.3 (44.62-3.60) 98.5 (49.36-3.60)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.23 (at 3.57Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158, PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.233 , 0.270 0.233 , 0.268	Depositor DCC
R_{free} test set	2001 reflections (3.07%)	wwPDB-VP
Wilson B-factor (Å ²)	115.3	Xtriage
Anisotropy	0.514	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 91.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	28400	wwPDB-VP
Average B, all atoms (Å ²)	144.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.44% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.25	0/1814	0.51	0/2466
1	B	0.25	0/1782	0.52	0/2424
2	C	0.25	0/8937	0.52	0/12087
3	D	0.25	0/11944	0.50	0/16149
4	E	0.24	0/775	0.47	0/1045
5	F	0.24	0/2852	0.49	0/3837
6	G	0.54	0/340	0.92	0/522
7	H	0.50	0/423	0.87	0/651
8	I	0.31	0/115	0.92	0/176
All	All	0.26	0/28982	0.53	0/39357

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1782	0	1834	27	0
1	B	1750	0	1797	19	0
2	C	8770	0	8874	113	0
3	D	11738	0	11971	146	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	E	761	0	778	14	0
5	F	2807	0	2882	36	0
6	G	305	0	170	11	0
7	H	376	0	205	14	0
8	I	104	0	54	7	0
9	B	1	0	0	0	0
9	D	3	0	0	0	0
9	F	1	0	0	0	0
10	D	2	0	0	0	0
All	All	28400	0	28565	338	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 6.

All (338) 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)
6:G:15:DT:H2'	6:G:16:DG:C8	2.18	0.78
1:B:206:THR:HG22	1:B:209:GLU:H	1.50	0.76
2:C:1116:ALA:HB2	3:D:88:TYR:HB3	1.69	0.74
3:D:1491:THR:HG21	4:E:89:MET:HG2	1.73	0.71
2:C:628:PHE:H	2:C:638:ASP:HB2	1.55	0.70
3:D:238:PRO:HD3	3:D:318:ARG:HG3	1.73	0.70
2:C:905:ILE:HG23	2:C:906:PHE:HD2	1.54	0.70
2:C:462:ASP:HB3	2:C:468:ARG:HD2	1.75	0.69
1:A:94:LEU:O	1:A:146:ARG:NH1	2.26	0.68
2:C:55:GLU:O	2:C:56:GLU:HB3	1.94	0.68
3:D:61:GLY:O	3:D:64:LYS:NZ	2.27	0.68
2:C:983:ILE:HG21	2:C:987:ILE:HD11	1.76	0.67
3:D:124:GLU:OE2	3:D:587:ARG:NH1	2.27	0.67
3:D:1254:GLN:HB3	3:D:1258:ARG:HB2	1.77	0.67
6:G:11:DT:H2'	6:G:12:DC:C6	2.30	0.66
3:D:266:GLU:HG3	3:D:314:PRO:HB3	1.77	0.66
5:F:365:GLU:HB2	5:F:404:ALA:HB2	1.78	0.66
5:F:397:ILE:HD12	5:F:400:ILE:HD11	1.79	0.65
2:C:766:GLU:HG3	3:D:64:LYS:HD2	1.78	0.64
3:D:1495:ILE:HG12	4:E:88:GLU:HG3	1.79	0.64
3:D:65:ARG:NH1	5:F:378:GLY:O	2.28	0.64
1:A:222:LEU:HD21	1:B:218:LEU:HD23	1.80	0.64
3:D:520:LEU:O	3:D:525:ARG:NH1	2.31	0.64
1:A:179:PHE:HB3	1:A:197:LEU:HD23	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:371:ILE:HG23	5:F:230:LYS:HD2	1.79	0.64
3:D:1108:ARG:NH1	3:D:1198:TYR:O	2.31	0.64
7:H:20:DC:H2'	7:H:21:DA:C8	2.33	0.64
2:C:711:GLU:HG2	2:C:822:VAL:HG22	1.80	0.63
5:F:400:ILE:HA	5:F:403:LYS:HG2	1.80	0.63
2:C:343:GLN:HG3	2:C:385:PHE:HB2	1.81	0.62
7:H:17:DA:H2'	7:H:18:DT:C6	2.34	0.62
2:C:17:PRO:HB2	2:C:20:GLU:HB3	1.81	0.62
3:D:260:GLU:OE1	3:D:273:ARG:NH1	2.28	0.62
2:C:521:PRO:HB3	3:D:1068:LEU:HD21	1.81	0.61
3:D:643:GLY:HA3	3:D:727:GLN:HB2	1.83	0.61
3:D:1236:LEU:HA	3:D:1359:GLN:HG3	1.82	0.61
2:C:97:ARG:NH1	2:C:110:GLU:OE2	2.34	0.60
3:D:218:LYS:HG2	3:D:338:GLU:HG2	1.81	0.60
2:C:591:SER:O	2:C:593:ALA:N	2.26	0.60
2:C:936:VAL:HG11	2:C:959:PRO:HB2	1.83	0.60
2:C:1092:LEU:HD13	2:C:1099:VAL:HG21	1.84	0.60
1:A:24:VAL:HG22	1:A:196:THR:HG23	1.82	0.60
2:C:420:ARG:HH22	8:I:1:A:H5''	1.67	0.59
3:D:1101:VAL:HG13	3:D:1102:THR:HG23	1.85	0.59
1:A:106:PRO:HG3	1:A:134:GLU:HG2	1.83	0.59
4:E:95:VAL:HG12	4:E:95:VAL:O	2.01	0.59
2:C:428:ARG:NH2	2:C:447:ALA:O	2.34	0.59
2:C:397:GLU:HG3	2:C:631:SER:HB2	1.84	0.59
5:F:131:VAL:HG13	5:F:178:ARG:HD3	1.85	0.59
5:F:361:LEU:HB3	5:F:365:GLU:HG3	1.85	0.58
2:C:504:GLU:HG2	2:C:509:ALA:HB2	1.86	0.58
3:D:142:LEU:HB2	3:D:161:LEU:HD11	1.86	0.58
5:F:323:ASP:HB2	8:I:1:A:O4'	2.04	0.58
2:C:787:ASP:OD2	2:C:791:ARG:NH2	2.36	0.58
3:D:657:LEU:HG	3:D:661:MET:HE2	1.85	0.58
2:C:210:GLU:HB3	2:C:211:LEU:HD12	1.85	0.58
2:C:413:LEU:HD21	2:C:451:LEU:HD13	1.85	0.57
7:H:21:DA:H2''	7:H:22:DG:H5''	1.85	0.57
1:B:71:VAL:HG22	1:B:132:LEU:HG	1.87	0.57
3:D:480:GLU:OE2	3:D:488:ARG:NH1	2.37	0.57
6:G:15:DT:H2'	6:G:16:DG:H8	1.66	0.57
2:C:617:ASP:OD1	2:C:617:ASP:N	2.37	0.56
3:D:1353:GLN:NE2	3:D:1365:ASP:OD1	2.35	0.56
7:H:18:DT:H2''	7:H:19:DG:H8	1.69	0.56
2:C:92:ALA:HB2	2:C:120:LEU:HD11	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:F:153:PRO:HA	5:F:156:VAL:HG22	1.86	0.56
3:D:1461:GLY:O	3:D:1465:ASN:ND2	2.38	0.56
5:F:187:LEU:HD23	5:F:224:VAL:HG13	1.87	0.56
5:F:188:ILE:HD13	5:F:221:ILE:HG12	1.88	0.56
3:D:134:VAL:HG12	3:D:454:ALA:HB2	1.88	0.56
1:B:216:GLU:OE2	1:B:219:ARG:NH2	2.35	0.55
3:D:5:VAL:O	3:D:1470:ARG:NH2	2.40	0.54
3:D:1379:VAL:HG21	3:D:1400:VAL:HG11	1.89	0.54
3:D:904:VAL:HG22	3:D:905:PRO:HD2	1.90	0.54
3:D:959:GLU:N	3:D:959:GLU:OE1	2.40	0.54
2:C:259:GLY:HA2	2:C:263:ASP:HB2	1.90	0.54
2:C:773:LEU:HB2	5:F:375:LEU:HD11	1.90	0.54
3:D:832:ARG:HD2	3:D:833:GLU:H	1.72	0.54
2:C:605:LYS:HB2	2:C:612:VAL:HB	1.90	0.54
3:D:841:TYR:HB2	3:D:864:VAL:HG22	1.89	0.53
3:D:462:GLN:HB2	3:D:513:ILE:HG21	1.90	0.53
3:D:1258:ARG:HH21	3:D:1351:GLU:HG2	1.72	0.53
2:C:557:ARG:HG3	2:C:844:GLY:HA3	1.91	0.53
3:D:939:PHE:O	3:D:943:THR:HG22	2.09	0.53
2:C:168:ARG:NH2	2:C:265:ARG:O	2.42	0.53
5:F:200:LYS:HD3	7:H:8:DG:H5''	1.91	0.53
2:C:177:GLU:HG3	2:C:178:PRO:HD2	1.89	0.53
2:C:221:LEU:HD11	2:C:307:LEU:HD21	1.90	0.53
2:C:607:ASP:HB2	2:C:610:ARG:NH1	2.24	0.52
1:B:91:ASN:HB3	1:B:94:LEU:HB2	1.91	0.52
2:C:1056:LYS:O	3:D:624:ASP:N	2.38	0.52
3:D:1042:ARG:HB3	3:D:1057:VAL:HB	1.92	0.52
2:C:420:ARG:NH1	8:I:1:A:OP1	2.42	0.52
2:C:709:GLU:OE2	2:C:824:ARG:NH1	2.43	0.52
2:C:418:LEU:HD21	2:C:427:VAL:HG11	1.91	0.52
1:A:198:ARG:HD3	2:C:934:PHE:CZ	2.45	0.52
2:C:1019:GLN:HG2	2:C:1058:ASP:HB3	1.91	0.52
3:D:242:LEU:HB3	3:D:311:LEU:HD12	1.92	0.52
2:C:203:ASP:OD1	2:C:204:GLN:N	2.43	0.51
3:D:622:ARG:NH1	6:G:17:DA:OP1	2.38	0.51
7:H:20:DC:H2'	7:H:21:DA:H8	1.76	0.51
1:B:220:GLU:O	1:B:223:THR:OG1	2.22	0.51
2:C:1037:VAL:HG13	2:C:1049:LEU:HD11	1.91	0.51
1:B:56:VAL:HG21	1:B:82:LEU:HD13	1.92	0.51
5:F:393:THR:HG22	5:F:395:GLU:H	1.76	0.50
1:A:206:THR:HG22	1:A:208:LEU:H	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:H:19:DG:H2''	7:H:20:DC:H6	1.76	0.50
2:C:774:LEU:HD23	5:F:354:LEU:HD21	1.92	0.50
3:D:42:ASP:OD1	3:D:48:ARG:NH2	2.44	0.50
3:D:1046:GLN:OE1	3:D:1046:GLN:N	2.43	0.50
3:D:561:GLY:HA2	5:F:140:ARG:HH11	1.75	0.50
3:D:1068:LEU:O	3:D:1072:ILE:HG12	2.12	0.50
2:C:194:VAL:HG22	2:C:221:LEU:HD23	1.93	0.50
3:D:1100:ASP:OD1	3:D:1463:LYS:NZ	2.31	0.50
2:C:593:ALA:HB1	2:C:659:PRO:HD2	1.94	0.50
2:C:876:VAL:HB	3:D:949:ILE:HD12	1.94	0.49
3:D:356:PRO:HG2	3:D:359:ALA:HB2	1.94	0.49
3:D:350:HIS:HE1	5:F:96:LEU:HD11	1.78	0.49
2:C:802:ARG:HB2	2:C:826:TYR:HB2	1.93	0.49
3:D:41:ARG:HG3	3:D:48:ARG:HE	1.76	0.49
3:D:520:LEU:HD23	3:D:525:ARG:HG2	1.93	0.49
3:D:474:GLU:HG3	3:D:496:LEU:HD11	1.95	0.49
3:D:1150:ALA:HB3	3:D:1187:PRO:HB2	1.95	0.49
3:D:1493:LYS:O	3:D:1497:GLU:HG2	2.13	0.49
6:G:14:DC:H2'	6:G:15:DT:C6	2.48	0.49
1:A:53:VAL:HG22	1:A:144:VAL:HG22	1.95	0.49
3:D:171:LEU:HD22	3:D:390:PRO:HG2	1.95	0.49
3:D:760:ARG:NH1	4:E:59:ASN:OD1	2.45	0.49
1:A:20:TYR:OH	1:A:198:ARG:HD2	2.13	0.48
2:C:578:VAL:HG23	2:C:579:VAL:HG23	1.95	0.48
3:D:809:PRO:HB3	3:D:839:LEU:HD13	1.95	0.48
1:B:56:VAL:HG22	1:B:142:VAL:HG12	1.95	0.48
3:D:897:TRP:CH2	3:D:902:LEU:HD22	2.48	0.48
2:C:740:GLU:HB3	2:C:805:ARG:HH12	1.78	0.48
1:A:32:PHE:HA	1:A:35:THR:HB	1.96	0.48
1:B:32:PHE:HA	1:B:35:THR:HB	1.96	0.48
2:C:1104:GLU:HB3	3:D:6:ARG:HG3	1.96	0.48
5:F:400:ILE:HG22	5:F:403:LYS:HE2	1.96	0.48
5:F:163:LEU:HD13	5:F:174:LEU:HD13	1.95	0.48
3:D:1102:THR:HG21	3:D:1371:VAL:HG22	1.96	0.48
6:G:19:DT:H3	8:I:1:A:H61	1.62	0.48
2:C:772:ARG:NE	5:F:380:GLU:OE2	2.47	0.47
2:C:577:PRO:HG2	2:C:580:MET:HG2	1.95	0.47
3:D:233:LYS:NZ	3:D:240:GLU:OE2	2.31	0.47
2:C:168:ARG:HD3	2:C:268:ASP:HB3	1.95	0.47
3:D:996:TRP:CD2	3:D:1056:PRO:HG3	2.48	0.47
1:B:49:PRO:HA	1:B:148:VAL:HG12	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:216:GLU:OE2	1:A:219:ARG:NH1	2.47	0.47
2:C:200:LEU:HD13	2:C:300:ASP:HB2	1.96	0.47
3:D:475:LYS:O	3:D:479:GLU:HG2	2.14	0.47
2:C:915:LYS:NZ	3:D:952:ASP:OD2	2.46	0.47
3:D:1000:THR:HG23	3:D:1036:ARG:HD2	1.97	0.47
2:C:627:ARG:NE	2:C:639:GLN:O	2.46	0.47
2:C:942:GLU:HG2	2:C:945:ARG:HH21	1.79	0.47
2:C:1067:TYR:OH	3:D:674:ARG:NH1	2.48	0.47
3:D:1137:ARG:O	3:D:1141:GLU:HG3	2.15	0.47
5:F:270:LYS:HG2	5:F:295:MET:HE1	1.96	0.47
1:B:124:ASN:OD1	1:B:124:ASN:N	2.48	0.47
3:D:1381:VAL:HG21	3:D:1389:LEU:HD23	1.96	0.47
2:C:211:LEU:HD11	2:C:304:LEU:HD11	1.97	0.47
1:A:64:GLU:HG2	1:A:76:VAL:HG22	1.97	0.46
7:H:18:DT:H2"	7:H:19:DG:C8	2.49	0.46
2:C:704:HIS:CD2	2:C:831:ARG:HD2	2.50	0.46
2:C:911:GLU:O	2:C:915:LYS:HG2	2.16	0.46
3:D:236:TYR:HD2	3:D:322:VAL:HG21	1.79	0.46
3:D:90:MET:SD	3:D:521:PRO:HD3	2.56	0.46
4:E:83:ASP:OD1	4:E:83:ASP:N	2.47	0.46
2:C:1070:ILE:CG2	3:D:655:PRO:HB2	2.46	0.46
3:D:455:ARG:HB2	3:D:460:ALA:HB2	1.97	0.46
3:D:689:ASP:O	3:D:693:GLU:HG3	2.16	0.46
1:A:100:LEU:HD23	1:A:141:GLU:HG2	1.98	0.46
5:F:166:LEU:HD13	5:F:170:HIS:HB3	1.97	0.46
2:C:1115:LEU:HG	3:D:85:VAL:HG12	1.98	0.46
3:D:244:GLU:HG3	3:D:310:LEU:HG	1.96	0.46
3:D:619:LEU:HD11	3:D:1439:SER:HB2	1.98	0.46
3:D:703:ASN:HB2	3:D:713:ILE:HG12	1.97	0.46
3:D:715:ALA:HB3	3:D:764:LEU:HA	1.97	0.46
3:D:829:VAL:HG21	3:D:839:LEU:HD11	1.98	0.46
5:F:397:ILE:HA	5:F:400:ILE:HG12	1.96	0.46
2:C:395:LYS:HE2	2:C:403:SER:HB3	1.98	0.46
2:C:1031:ARG:HG2	6:G:16:DG:H5"	1.97	0.46
5:F:324:GLU:HG3	8:I:1:A:C8	2.51	0.46
1:A:107:LYS:HE3	1:A:107:LYS:HB2	1.80	0.45
1:A:41:ARG:HA	1:A:177:VAL:HG11	1.98	0.45
3:D:821:VAL:HG11	3:D:827:ILE:HD12	1.98	0.45
6:G:6:DC:C6	6:G:7:DT:H72	2.52	0.45
3:D:273:ARG:HB3	3:D:278:PRO:HA	1.99	0.45
3:D:1485:GLN:NE2	4:E:82:GLU:OE1	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:9:ARG:HB2	3:D:1456:LYS:HG2	1.99	0.45
3:D:1084:THR:O	3:D:1088:THR:HG23	2.17	0.45
5:F:80:PRO:HB2	5:F:210:LEU:HD11	1.99	0.45
2:C:81:ASP:OD1	2:C:81:ASP:N	2.39	0.45
2:C:719:PRO:HB3	2:C:820:ARG:NE	2.31	0.45
5:F:236:SER:OG	7:H:5:DA:OP2	2.34	0.45
1:B:108:GLU:HG2	1:B:131:THR:HG22	1.99	0.45
2:C:164:PRO:HD2	2:C:171:TRP:CD1	2.51	0.45
2:C:165:LEU:HB2	2:C:168:ARG:HG3	1.98	0.45
4:E:14:ASP:OD1	4:E:14:ASP:N	2.50	0.45
1:A:180:GLN:NE2	2:C:935:GLY:O	2.49	0.44
1:A:10:VAL:HG22	1:A:26:GLU:O	2.17	0.44
1:B:175:ARG:N	1:B:200:TRP:O	2.44	0.44
3:D:17:LYS:HB2	3:D:17:LYS:HE3	1.74	0.44
3:D:236:TYR:CD2	3:D:322:VAL:HG21	2.51	0.44
5:F:164:LYS:HA	5:F:171:LYS:HE3	2.00	0.44
2:C:740:GLU:HB3	2:C:805:ARG:NH1	2.32	0.44
3:D:618:LEU:HG	3:D:1467:ILE:HG23	2.00	0.44
1:B:92:PRO:O	1:B:146:ARG:NH1	2.44	0.44
2:C:212:GLY:HA2	2:C:218:VAL:HG21	1.99	0.44
3:D:101:HIS:HB3	3:D:104:PHE:HD2	1.82	0.44
3:D:834:THR:OG1	3:D:835:SER:N	2.48	0.44
2:C:922:PHE:CE2	2:C:964:LYS:HB2	2.53	0.44
2:C:1009:SER:HB3	3:D:651:GLU:O	2.16	0.44
4:E:33:HIS:CE1	4:E:89:MET:HB3	2.53	0.44
5:F:330:GLY:HA2	5:F:333:ILE:HD12	2.00	0.44
3:D:1147:ARG:NH2	3:D:1369:GLU:OE1	2.48	0.44
3:D:34:TYR:CZ	3:D:35:ARG:HG3	2.53	0.44
3:D:355:VAL:HG11	3:D:385:VAL:HG21	1.99	0.44
1:B:80:LEU:HG	3:D:844:ALA:HA	1.99	0.44
3:D:191:LEU:HB3	3:D:393:ILE:HD12	2.00	0.44
3:D:477:LEU:HB2	3:D:496:LEU:HD13	1.99	0.44
3:D:1387:SER:HB3	3:D:1407:LEU:HD11	2.00	0.44
3:D:1098:LEU:HA	3:D:1101:VAL:HG12	2.00	0.43
3:D:355:VAL:HG13	3:D:359:ALA:HB3	2.00	0.43
3:D:850:LEU:HD12	3:D:884:ARG:NH2	2.34	0.43
4:E:66:LYS:O	4:E:70:THR:HG23	2.18	0.43
7:H:16:DG:H2"	7:H:17:DA:C8	2.53	0.43
1:A:8:ALA:HA	1:A:9:PRO:HD3	1.84	0.43
2:C:247:PRO:HA	2:C:248:PRO:HD3	1.75	0.43
2:C:598:GLU:O	2:C:651:LYS:HG3	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:1377:LYS:HE3	3:D:1378:TYR:CZ	2.54	0.43
3:D:236:TYR:CZ	3:D:242:LEU:HD12	2.54	0.43
3:D:1372:VAL:HA	3:D:1375:MET:HE3	2.00	0.43
1:B:153:ALA:HA	1:B:156:HIS:NE2	2.33	0.43
2:C:64:LEU:HD22	2:C:100:LEU:HD11	1.99	0.43
3:D:248:PRO:HG3	3:D:308:LYS:HE3	2.00	0.43
3:D:1495:ILE:HD13	4:E:80:VAL:HG21	2.00	0.43
3:D:684:LYS:HE2	3:D:684:LYS:HB3	1.87	0.43
3:D:1053:PHE:CZ	3:D:1072:ILE:HD12	2.53	0.43
2:C:160:ALA:HB3	2:C:174:LEU:HB2	1.99	0.43
2:C:229:MET:HB2	2:C:233:GLU:HB2	2.01	0.43
3:D:14:SER:HB3	3:D:511:TRP:CE2	2.54	0.43
3:D:462:GLN:NE2	3:D:515:GLU:OE2	2.46	0.43
3:D:960:LYS:NZ	3:D:1063:GLU:OE2	2.46	0.43
5:F:278:LEU:HA	5:F:281:GLU:HG2	1.99	0.43
1:B:64:GLU:HA	1:B:165:ILE:HD13	2.00	0.43
3:D:26:VAL:HG11	3:D:44:LEU:HD23	2.01	0.43
2:C:15:LEU:O	2:C:586:ARG:NH2	2.45	0.43
3:D:230:TRP:CZ2	3:D:232:GLU:HG2	2.54	0.43
4:E:52:GLU:OE1	4:E:52:GLU:N	2.46	0.43
5:F:237:THR:OG1	7:H:4:DA:H2'	2.18	0.43
2:C:535:SER:O	2:C:538:GLN:HG2	2.19	0.42
1:A:133:GLU:HG2	1:A:134:GLU:N	2.34	0.42
2:C:49:ARG:CZ	2:C:49:ARG:HB3	2.46	0.42
2:C:577:PRO:HB3	2:C:993:PHE:CG	2.54	0.42
2:C:1043:TYR:CG	3:D:763:MET:HG2	2.53	0.42
3:D:192:ALA:HB3	3:D:195:VAL:HB	2.00	0.42
3:D:787:LEU:HD21	3:D:947:ILE:HG21	2.00	0.42
4:E:46:PRO:HD2	4:E:63:TRP:CE2	2.54	0.42
3:D:314:PRO:HB2	3:D:317:VAL:HG12	2.00	0.42
3:D:784:ASP:HB2	3:D:939:PHE:HE1	1.84	0.42
3:D:1277:ILE:HG22	3:D:1278:ASP:H	1.84	0.42
1:A:206:THR:HB	1:A:209:GLU:HG3	2.02	0.42
2:C:843:HIS:NE2	2:C:887:GLU:OE1	2.53	0.42
1:A:176:ARG:NH1	2:C:863:ASP:O	2.52	0.42
3:D:1208:ASP:HB2	3:D:1215:VAL:HA	2.01	0.42
5:F:323:ASP:OD1	8:I:1:A:H1'	2.19	0.42
2:C:684:PHE:HB3	3:D:633:VAL:HG21	2.00	0.42
3:D:983:LEU:HD13	3:D:988:ARG:HB2	2.01	0.42
3:D:1463:LYS:HE2	3:D:1463:LYS:HB3	1.86	0.42
3:D:661:MET:HE3	3:D:673:ALA:HB1	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:784:ASP:HB2	3:D:939:PHE:CE1	2.55	0.42
1:B:111:ALA:HB3	1:B:125:PRO:HA	2.02	0.42
2:C:11:GLU:H	2:C:11:GLU:HG2	1.50	0.42
2:C:12:VAL:HG21	2:C:472:ARG:HD3	2.02	0.42
2:C:351:LEU:HD12	2:C:375:SER:HA	2.01	0.42
3:D:808:THR:HB	3:D:810:GLU:HG2	2.01	0.42
2:C:516:ARG:HD3	3:D:1068:LEU:HD13	2.02	0.42
2:C:1066:ALA:O	2:C:1070:ILE:HG13	2.20	0.42
2:C:258:TYR:O	2:C:263:ASP:N	2.53	0.42
3:D:889:ALA:HB1	3:D:930:LEU:HA	2.02	0.42
1:A:57:TYR:CG	1:A:161:ARG:HD2	2.56	0.41
3:D:44:LEU:HB3	3:D:525:ARG:NH2	2.35	0.41
3:D:131:LYS:O	3:D:456:MET:HG2	2.19	0.41
3:D:1126:ASP:OD1	3:D:1128:VAL:HG13	2.20	0.41
2:C:797:GLY:O	2:C:829:GLN:NE2	2.53	0.41
3:D:615:ARG:HH12	6:G:13:DC:P	2.43	0.41
3:D:1040:GLY:O	3:D:1060:SER:HB3	2.20	0.41
3:D:1491:THR:O	3:D:1495:ILE:HG13	2.20	0.41
7:H:19:DG:H2''	7:H:20:DC:C6	2.54	0.41
2:C:103:LYS:H	2:C:103:LYS:HG3	1.60	0.41
2:C:304:LEU:HB3	2:C:305:PRO:HD3	2.02	0.41
2:C:351:LEU:HD11	2:C:373:VAL:HG13	2.02	0.41
3:D:411:THR:HB	3:D:437:VAL:H	1.85	0.41
8:I:1:A:H2'	8:I:2:C:C6	2.55	0.41
1:A:104:GLU:OE1	1:A:137:ARG:NH1	2.53	0.41
2:C:808:ARG:NH1	5:F:305:GLU:OE2	2.53	0.41
3:D:1189:ARG:HB3	3:D:1204:CYS:HA	2.02	0.41
3:D:1252:ILE:HG23	3:D:1253:THR:HG23	2.02	0.41
4:E:40:LEU:HG	4:E:67:GLU:HG2	2.01	0.41
5:F:172:ARG:O	5:F:176:ILE:HG12	2.20	0.41
1:A:36:LEU:HD23	1:A:36:LEU:HA	1.92	0.41
3:D:114:THR:HG23	3:D:495:ARG:HG2	2.02	0.41
2:C:545:ASN:HB3	2:C:583:LEU:HD22	2.02	0.41
3:D:1047:LYS:HG2	3:D:1053:PHE:CZ	2.55	0.41
2:C:107:LEU:HD22	2:C:108:ILE:N	2.36	0.41
2:C:324:ASP:HB3	2:C:327:HIS:HB2	2.02	0.41
2:C:503:LEU:HD23	2:C:508:ILE:HA	2.02	0.41
3:D:828:LYS:HG2	3:D:833:GLU:HB3	2.03	0.41
6:G:7:DT:H2''	6:G:8:DG:H8	1.86	0.41
2:C:146:VAL:HG11	2:C:306:THR:HG22	2.01	0.41
2:C:348:LEU:HD12	2:C:348:LEU:HA	1.96	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:886:LEU:HD21	3:D:951:ILE:HG12	2.02	0.41
3:D:438:ASP:OD2	3:D:441:ARG:NH1	2.50	0.41
3:D:895:VAL:O	3:D:899:LEU:HG	2.21	0.41
3:D:1053:PHE:CE1	3:D:1072:ILE:HG23	2.56	0.41
1:A:70:GLY:N	2:C:607:ASP:OD1	2.53	0.41
1:B:58:ILE:HB	1:B:61:VAL:HB	2.03	0.41
2:C:133:ASP:HB3	2:C:395:LYS:HD2	2.03	0.41
2:C:136:ILE:HB	2:C:336:VAL:HG13	2.03	0.41
2:C:430:VAL:O	3:D:1075:HIS:ND1	2.50	0.41
2:C:501:THR:HA	2:C:502:PRO:HD3	1.95	0.41
2:C:605:LYS:HB3	2:C:610:ARG:NH1	2.35	0.41
2:C:905:ILE:HG23	2:C:906:PHE:CD2	2.44	0.41
3:D:654:LYS:O	3:D:658:LEU:HG	2.20	0.41
3:D:1152:GLU:HG3	3:D:1161:GLU:HA	2.03	0.41
3:D:1495:ILE:HG22	3:D:1499:ARG:HD2	2.02	0.41
4:E:47:LYS:HA	4:E:47:LYS:HD3	1.90	0.41
5:F:188:ILE:HG12	5:F:224:VAL:HG21	2.02	0.41
1:A:153:ALA:HA	1:A:156:HIS:NE2	2.35	0.41
1:A:57:TYR:CD1	1:A:161:ARG:HD2	2.57	0.40
6:G:11:DT:H2"	6:G:12:DC:OP1	2.21	0.40
2:C:896:PHE:HB2	2:C:921:ALA:HB1	2.04	0.40
3:D:1099:VAL:O	3:D:1103:HIS:HB3	2.21	0.40
3:D:553:ARG:HD3	5:F:214:GLN:HB3	2.04	0.40
2:C:35:PRO:HA	2:C:36:PRO:HD3	1.96	0.40
2:C:218:VAL:O	2:C:222:MET:HG2	2.21	0.40
7:H:7:DG:H5'	7:H:7:DG:H8	1.86	0.40
5:F:193:ARG:HB3	7:H:7:DG:H5"	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	224/315 (71%)	222 (99%)	2 (1%)	0	100	100
1	B	220/315 (70%)	215 (98%)	5 (2%)	0	100	100
2	C	1107/1119 (99%)	1088 (98%)	18 (2%)	1 (0%)	51	83
3	D	1482/1524 (97%)	1459 (98%)	21 (1%)	2 (0%)	51	83
4	E	92/99 (93%)	91 (99%)	1 (1%)	0	100	100
5	F	344/443 (78%)	340 (99%)	4 (1%)	0	100	100
All	All	3469/3815 (91%)	3415 (98%)	51 (2%)	3 (0%)	51	83

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	592	LEU
3	D	1440	PHE
3	D	530	VAL

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	199/273 (73%)	192 (96%)	7 (4%)	36	68
1	B	195/273 (71%)	190 (97%)	5 (3%)	46	74
2	C	936/941 (100%)	877 (94%)	59 (6%)	18	53
3	D	1253/1279 (98%)	1179 (94%)	74 (6%)	19	55
4	E	83/88 (94%)	82 (99%)	1 (1%)	71	87
5	F	301/388 (78%)	284 (94%)	17 (6%)	21	56
All	All	2967/3242 (92%)	2804 (94%)	163 (6%)	21	57

All (163) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LEU
1	A	34	VAL

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Mol	Chain	Res	Type
1	A	67	THR
1	A	142	VAL
1	A	193	ASP
1	A	219	ARG
1	A	229	GLN
1	B	34	VAL
1	B	80	LEU
1	B	94	LEU
1	B	186	LEU
1	B	206	THR
2	C	8	ARG
2	C	11	GLU
2	C	15	LEU
2	C	49	ARG
2	C	56	GLU
2	C	81	ASP
2	C	97	ARG
2	C	103	LYS
2	C	107	LEU
2	C	133	ASP
2	C	141	HIS
2	C	168	ARG
2	C	177	GLU
2	C	205	GLU
2	C	221	LEU
2	C	232	GLU
2	C	251	ASP
2	C	274	ARG
2	C	331	ARG
2	C	342	ASP
2	C	358	ARG
2	C	403	SER
2	C	418	LEU
2	C	419	THR
2	C	427	VAL
2	C	434	HIS
2	C	464	LEU
2	C	480	THR
2	C	482	GLU
2	C	512	ARG
2	C	524	VAL
2	C	557	ARG

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Mol	Chain	Res	Type
2	C	575	GLN
2	C	583	LEU
2	C	610	ARG
2	C	617	ASP
2	C	633	GLN
2	C	638	ASP
2	C	640	ARG
2	C	648	ARG
2	C	680	ASP
2	C	774	LEU
2	C	775	ARG
2	C	786	LYS
2	C	807	ARG
2	C	808	ARG
2	C	813	VAL
2	C	815	LEU
2	C	820	ARG
2	C	830	LYS
2	C	928	LYS
2	C	939	ARG
2	C	952	LEU
2	C	968	LEU
2	C	978	ARG
2	C	1001	VAL
2	C	1057	SER
2	C	1058	ASP
2	C	1115	LEU
3	D	67	ARG
3	D	68	PHE
3	D	80	VAL
3	D	81	THR
3	D	135	LEU
3	D	141	ILE
3	D	155	ASP
3	D	161	LEU
3	D	190	GLU
3	D	191	LEU
3	D	198	ARG
3	D	200	ASP
3	D	230	TRP
3	D	231	VAL
3	D	256	GLU

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Mol	Chain	Res	Type
3	D	270	LEU
3	D	275	GLU
3	D	276	ASP
3	D	312	ARG
3	D	331	VAL
3	D	362	GLU
3	D	372	ASP
3	D	387	LEU
3	D	411	THR
3	D	421	LEU
3	D	430	ASP
3	D	486	ARG
3	D	500	ARG
3	D	525	ARG
3	D	572	ARG
3	D	576	GLU
3	D	586	ARG
3	D	591	VAL
3	D	618	LEU
3	D	632	VAL
3	D	650	LEU
3	D	669	ASN
3	D	709	HIS
3	D	754	PHE
3	D	778	LEU
3	D	808	THR
3	D	817	GLU
3	D	832	ARG
3	D	864	VAL
3	D	867	ARG
3	D	894	LYS
3	D	904	VAL
3	D	943	THR
3	D	983	LEU
3	D	984	THR
3	D	1041	LEU
3	D	1062	ARG
3	D	1127	GLU
3	D	1128	VAL
3	D	1130	ARG
3	D	1155	VAL
3	D	1162	GLU

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Mol	Chain	Res	Type
3	D	1188	VAL
3	D	1195	GLN
3	D	1219	GLU
3	D	1221	VAL
3	D	1277	ILE
3	D	1284	GLU
3	D	1287	GLU
3	D	1288	GLU
3	D	1290	LEU
3	D	1304	LYS
3	D	1305	LEU
3	D	1313	VAL
3	D	1317	ASP
3	D	1470	ARG
3	D	1493	LYS
3	D	1496	GLU
3	D	1501	GLU
4	E	50	THR
5	F	88	ILE
5	F	123	ASP
5	F	125	ASP
5	F	141	VAL
5	F	150	THR
5	F	172	ARG
5	F	186	HIS
5	F	205	ARG
5	F	218	GLN
5	F	279	GLN
5	F	287	THR
5	F	310	ILE
5	F	364	ARG
5	F	377	ASP
5	F	417	LYS
5	F	420	ASP
5	F	422	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
3	D	350	HIS
3	D	1195	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
8	I	4/5 (80%)	0	0

There are no RNA backbone outliers to report.

There are no RNA pucker outliers to report.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 7 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	226/315 (71%)	-0.09	1 (0%) 92 86	99, 144, 168, 177	0
1	B	222/315 (70%)	-0.18	1 (0%) 91 83	98, 143, 178, 192	0
2	C	1111/1119 (99%)	0.09	38 (3%) 45 30	87, 140, 186, 211	0
3	D	1486/1524 (97%)	0.17	71 (4%) 30 19	87, 136, 187, 216	1 (0%)
4	E	94/99 (94%)	-0.17	1 (1%) 80 68	97, 159, 197, 210	0
5	F	346/443 (78%)	0.17	17 (4%) 29 18	101, 155, 209, 223	0
6	G	15/21 (71%)	-0.22	0 100 100	116, 162, 249, 258	0
7	H	18/27 (66%)	-0.30	0 100 100	131, 162, 242, 256	0
8	I	5/5 (100%)	0.28	0 100 100	116, 117, 128, 147	0
All	All	3523/3868 (91%)	0.09	129 (3%) 41 27	87, 142, 189, 258	1 (0%)

All (129) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	D	322	VAL	6.7
2	C	415	PRO	6.0
3	D	335	LEU	5.8
3	D	262	LYS	5.8
3	D	337	LEU	5.6
3	D	368	VAL	5.5
3	D	282	TYR	5.5
2	C	100	LEU	4.8
3	D	219	GLU	4.8
3	D	269	PHE	4.7
2	C	66	LEU	4.7
3	D	367	ILE	4.4
3	D	353	VAL	4.3
3	D	409	VAL	4.2
3	D	1294	VAL	4.2

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Mol	Chain	Res	Type	RSRZ
3	D	355	VAL	4.0
5	F	327	SER	4.0
2	C	99	GLN	3.8
5	F	418	LEU	3.8
2	C	414	GLY	3.8
2	C	766	GLU	3.7
3	D	393	ILE	3.6
3	D	321	GLN	3.6
5	F	326	ASP	3.6
3	D	421	LEU	3.5
3	D	1312	LEU	3.5
3	D	1327	ARG	3.5
5	F	381	HIS	3.4
3	D	349	PRO	3.4
2	C	741	GLY	3.3
5	F	149	GLU	3.3
3	D	305	ALA	3.3
2	C	776	SER	3.2
2	C	68	PHE	3.2
2	C	372	LEU	3.2
3	D	350	HIS	3.1
3	D	220	ARG	3.1
2	C	54	ILE	3.1
5	F	321	ILE	3.1
3	D	993	LEU	3.0
3	D	1292	VAL	2.9
2	C	608	GLY	2.9
5	F	375	LEU	2.9
3	D	221	ALA	2.9
5	F	419	ARG	2.9
2	C	756	VAL	2.9
3	D	170	PRO	2.9
3	D	1318	TYR	2.8
3	D	64	LYS	2.8
2	C	174	LEU	2.8
3	D	153	LEU	2.8
3	D	276	ASP	2.7
2	C	611	ILE	2.7
3	D	223	LEU	2.7
1	B	64	GLU	2.7
3	D	445	ARG	2.6
2	C	207	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
5	F	421	PHE	2.6
3	D	152	LEU	2.6
3	D	406	ASP	2.6
2	C	755	LEU	2.6
5	F	320	PRO	2.6
3	D	1132	LEU	2.6
2	C	48	PHE	2.6
3	D	352	ASN	2.6
2	C	769	PRO	2.5
3	D	295	GLY	2.5
3	D	666	ILE	2.5
2	C	650	ARG	2.5
3	D	268	ALA	2.5
2	C	98	LEU	2.5
3	D	1313	VAL	2.5
5	F	97	GLU	2.4
3	D	687	VAL	2.4
2	C	52	PHE	2.4
2	C	643	VAL	2.4
2	C	773	LEU	2.4
3	D	334	THR	2.4
2	C	777	ILE	2.4
3	D	586	ARG	2.4
2	C	109	LYS	2.4
5	F	235	PHE	2.4
2	C	743	VAL	2.4
2	C	657	ASP	2.4
3	D	205	TYR	2.3
3	D	263	GLU	2.3
3	D	1300	SER	2.3
2	C	50	GLU	2.3
2	C	419	THR	2.3
3	D	211	VAL	2.2
3	D	379	ALA	2.2
1	A	160	ASP	2.2
3	D	304	LEU	2.2
2	C	71	TYR	2.2
2	C	649	VAL	2.2
3	D	280	ALA	2.2
3	D	422	ALA	2.2
3	D	446	VAL	2.2
5	F	323	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
3	D	191	LEU	2.2
3	D	283	PHE	2.2
3	D	1490	LYS	2.2
3	D	197	SER	2.2
5	F	404	ALA	2.2
2	C	67	ASP	2.2
3	D	1279	GLY	2.2
3	D	821	VAL	2.2
3	D	658	LEU	2.1
2	C	418	LEU	2.1
3	D	360	ARG	2.1
2	C	64	LEU	2.1
2	C	386	PHE	2.1
3	D	387	LEU	2.1
3	D	339	TRP	2.1
2	C	226	VAL	2.1
3	D	1488	ASP	2.1
3	D	452	ILE	2.1
3	D	1278	ASP	2.1
3	D	26	VAL	2.1
3	D	675	ARG	2.1
3	D	171	LEU	2.1
2	C	227	PHE	2.1
3	D	226	PRO	2.0
3	D	281	THR	2.0
4	E	79	LEU	2.0
3	D	390	PRO	2.0
5	F	163	LEU	2.0
5	F	413	SER	2.0
5	F	293	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
9	MG	B	2001	1/1	0.08	0.88	169,169,169,169	0
9	MG	D	2005	1/1	0.84	0.37	112,112,112,112	0
10	ZN	D	2002	1/1	0.90	0.08	215,215,215,215	0
9	MG	F	2001	1/1	0.91	0.06	132,132,132,132	0
9	MG	D	2003	1/1	0.93	0.33	86,86,86,86	0
9	MG	D	2004	1/1	0.94	0.55	86,86,86,86	0
10	ZN	D	2001	1/1	0.95	0.23	118,118,118,118	0

6.5 Other polymers [i](#)

There are no such residues in this entry.