



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:20 PM GMT

PDB ID : 3TNX
Title : Structure of the precursor of a thermostable variant of papain at 2.6 Angstroem resolution
Authors : Roy, S.; Choudhury, D.; Dattagupta, J.K.; Biswas, S.
Deposited on : 2011-09-02
Resolution : 2.62 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

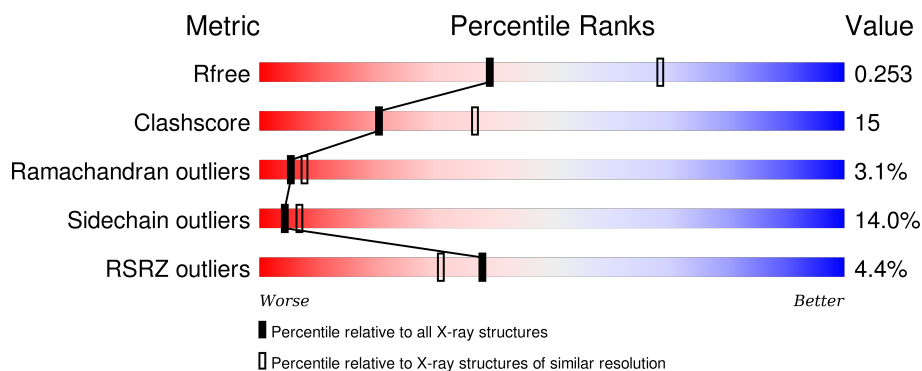
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.62 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	2700 (2.64-2.60)
Clashscore	102246	3065 (2.64-2.60)
Ramachandran outliers	100387	3015 (2.64-2.60)
Sidechain outliers	100360	3015 (2.64-2.60)
RSRZ outliers	91569	2706 (2.64-2.60)

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	363	 4% 60% 19% 6% 15%
1	C	363	 4% 53% 26% 6% 15%

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
2	CL	A	320	-	-	-	X

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5079 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 Papain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	310	Total	C	N	O	S	0	0	0
			2474	1567	423	476	8			
1	C	310	Total	C	N	O	S	0	0	0
			2471	1566	422	475	8			

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-43	MET	-	EXPRESSION TAG	UNP P00784
A	-42	HIS	-	EXPRESSION TAG	UNP P00784
A	-41	HIS	-	EXPRESSION TAG	UNP P00784
A	-40	HIS	-	EXPRESSION TAG	UNP P00784
A	-39	HIS	-	EXPRESSION TAG	UNP P00784
A	-38	HIS	-	EXPRESSION TAG	UNP P00784
A	-37	HIS	-	EXPRESSION TAG	UNP P00784
A	-36	SER	-	EXPRESSION TAG	UNP P00784
A	-35	SER	-	EXPRESSION TAG	UNP P00784
A	-34	GLY	-	EXPRESSION TAG	UNP P00784
A	-33	LEU	-	EXPRESSION TAG	UNP P00784
A	-32	VAL	-	EXPRESSION TAG	UNP P00784
A	-31	PRO	-	EXPRESSION TAG	UNP P00784
A	-30	ARG	-	EXPRESSION TAG	UNP P00784
A	-29	GLY	-	EXPRESSION TAG	UNP P00784
A	-28	SER	-	EXPRESSION TAG	UNP P00784
A	-27	GLY	-	EXPRESSION TAG	UNP P00784
A	-26	MET	-	EXPRESSION TAG	UNP P00784
A	-25	LYS	-	EXPRESSION TAG	UNP P00784
A	-24	GLU	-	EXPRESSION TAG	UNP P00784
A	-23	THR	-	EXPRESSION TAG	UNP P00784
A	-22	ALA	-	EXPRESSION TAG	UNP P00784
A	-21	ALA	-	EXPRESSION TAG	UNP P00784
A	-20	ALA	-	EXPRESSION TAG	UNP P00784
A	-19	LYS	-	EXPRESSION TAG	UNP P00784

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	-18	PHE	-	EXPRESSION TAG	UNP P00784
A	-17	GLU	-	EXPRESSION TAG	UNP P00784
A	-16	ARG	-	EXPRESSION TAG	UNP P00784
A	-15	GLN	-	EXPRESSION TAG	UNP P00784
A	-14	HIS	-	EXPRESSION TAG	UNP P00784
A	-13	MET	-	EXPRESSION TAG	UNP P00784
A	-12	ASP	-	EXPRESSION TAG	UNP P00784
A	-11	SER	-	EXPRESSION TAG	UNP P00784
A	-10	PRO	-	EXPRESSION TAG	UNP P00784
A	-9	ASP	-	EXPRESSION TAG	UNP P00784
A	-8	LEU	-	EXPRESSION TAG	UNP P00784
A	-7	GLY	-	EXPRESSION TAG	UNP P00784
A	-6	THR	-	EXPRESSION TAG	UNP P00784
A	-5	ASP	-	EXPRESSION TAG	UNP P00784
A	-4	ASP	-	EXPRESSION TAG	UNP P00784
A	-3	ASP	-	EXPRESSION TAG	UNP P00784
A	-2	ASP	-	EXPRESSION TAG	UNP P00784
A	-1	LYS	-	EXPRESSION TAG	UNP P00784
A	0	MET	-	EXPRESSION TAG	UNP P00784
A	132	ALA	CYS	ENGINEERED MUTATION	UNP P00784
A	139	SER	VAL	ENGINEERED MUTATION	UNP P00784
A	143	SER	GLY	ENGINEERED MUTATION	UNP P00784
A	281	ARG	LYS	ENGINEERED MUTATION	UNP P00784
C	-43	MET	-	EXPRESSION TAG	UNP P00784
C	-42	HIS	-	EXPRESSION TAG	UNP P00784
C	-41	HIS	-	EXPRESSION TAG	UNP P00784
C	-40	HIS	-	EXPRESSION TAG	UNP P00784
C	-39	HIS	-	EXPRESSION TAG	UNP P00784
C	-38	HIS	-	EXPRESSION TAG	UNP P00784
C	-37	HIS	-	EXPRESSION TAG	UNP P00784
C	-36	SER	-	EXPRESSION TAG	UNP P00784
C	-35	SER	-	EXPRESSION TAG	UNP P00784
C	-34	GLY	-	EXPRESSION TAG	UNP P00784
C	-33	LEU	-	EXPRESSION TAG	UNP P00784
C	-32	VAL	-	EXPRESSION TAG	UNP P00784
C	-31	PRO	-	EXPRESSION TAG	UNP P00784
C	-30	ARG	-	EXPRESSION TAG	UNP P00784
C	-29	GLY	-	EXPRESSION TAG	UNP P00784
C	-28	SER	-	EXPRESSION TAG	UNP P00784
C	-27	GLY	-	EXPRESSION TAG	UNP P00784
C	-26	MET	-	EXPRESSION TAG	UNP P00784
C	-25	LYS	-	EXPRESSION TAG	UNP P00784

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	-24	GLU	-	EXPRESSION TAG	UNP P00784
C	-23	THR	-	EXPRESSION TAG	UNP P00784
C	-22	ALA	-	EXPRESSION TAG	UNP P00784
C	-21	ALA	-	EXPRESSION TAG	UNP P00784
C	-20	ALA	-	EXPRESSION TAG	UNP P00784
C	-19	LYS	-	EXPRESSION TAG	UNP P00784
C	-18	PHE	-	EXPRESSION TAG	UNP P00784
C	-17	GLU	-	EXPRESSION TAG	UNP P00784
C	-16	ARG	-	EXPRESSION TAG	UNP P00784
C	-15	GLN	-	EXPRESSION TAG	UNP P00784
C	-14	HIS	-	EXPRESSION TAG	UNP P00784
C	-13	MET	-	EXPRESSION TAG	UNP P00784
C	-12	ASP	-	EXPRESSION TAG	UNP P00784
C	-11	SER	-	EXPRESSION TAG	UNP P00784
C	-10	PRO	-	EXPRESSION TAG	UNP P00784
C	-9	ASP	-	EXPRESSION TAG	UNP P00784
C	-8	LEU	-	EXPRESSION TAG	UNP P00784
C	-7	GLY	-	EXPRESSION TAG	UNP P00784
C	-6	THR	-	EXPRESSION TAG	UNP P00784
C	-5	ASP	-	EXPRESSION TAG	UNP P00784
C	-4	ASP	-	EXPRESSION TAG	UNP P00784
C	-3	ASP	-	EXPRESSION TAG	UNP P00784
C	-2	ASP	-	EXPRESSION TAG	UNP P00784
C	-1	LYS	-	EXPRESSION TAG	UNP P00784
C	0	MET	-	EXPRESSION TAG	UNP P00784
C	132	ALA	CYS	ENGINEERED MUTATION	UNP P00784
C	139	SER	VAL	ENGINEERED MUTATION	UNP P00784
C	143	SER	GLY	ENGINEERED MUTATION	UNP P00784
C	281	ARG	LYS	ENGINEERED MUTATION	UNP P00784

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Cl 1 1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	104	Total O 104 104	0	0

Continued on next page...

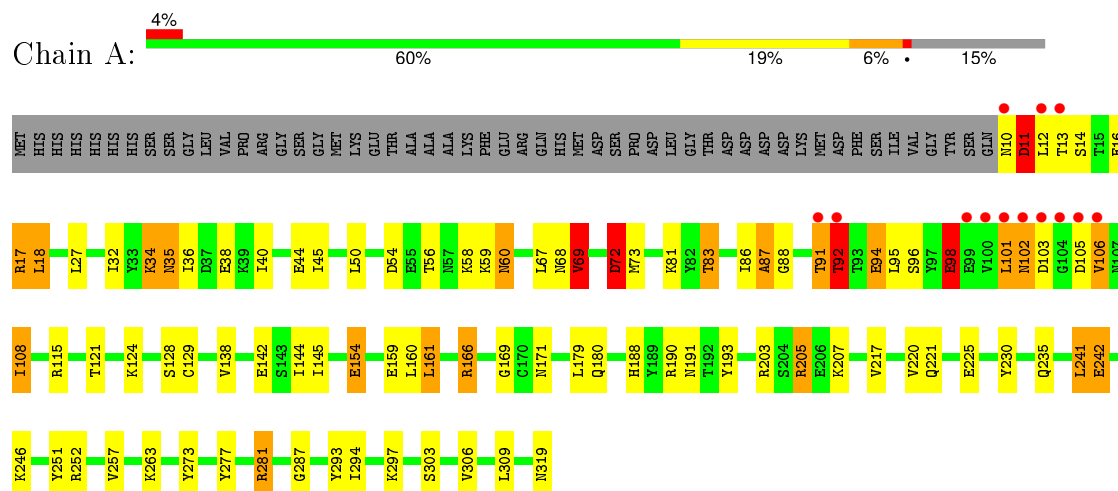
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	29	Total	O	0	0
			29	29		

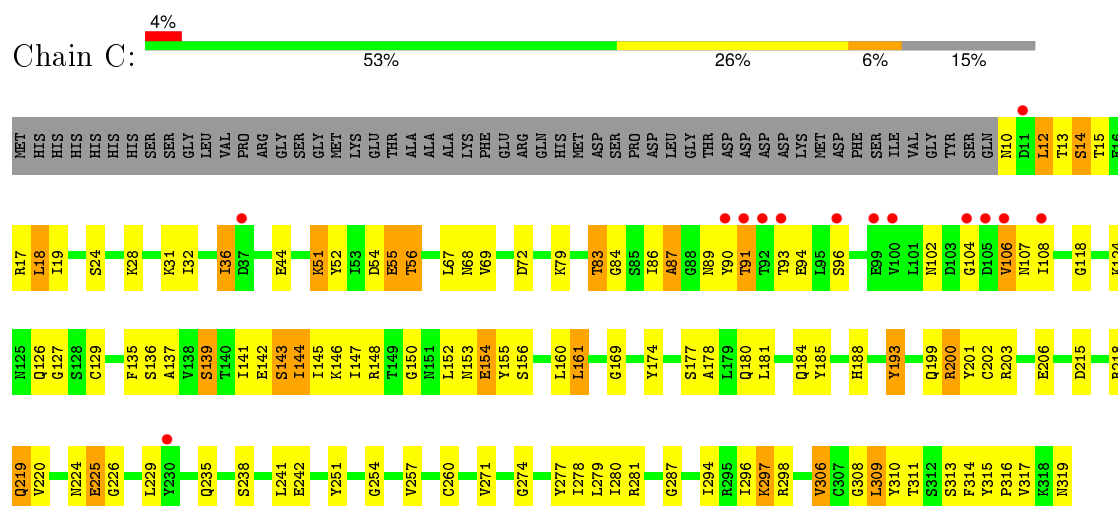
3 Residue-property plots

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: Papain



• Molecule 1: Papain



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	42.92Å 74.78Å 116.51Å 90.00° 93.03° 90.00°	Depositor
Resolution (Å)	45.92 – 2.62 45.91 – 2.62	Depositor EDS
% Data completeness (in resolution range)	96.5 (45.92-2.62) 96.5 (45.91-2.62)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.79 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.183 , 0.236 0.202 , 0.253	Depositor DCC
R_{free} test set	1106 reflections (5.39%)	DCC
Wilson B-factor (Å ²)	56.3	Xtriage
Anisotropy	0.454	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 57.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 21671 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5079	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.49% 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

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

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	0.95	4/2534 (0.2%)	0.87	5/3433 (0.1%)
1	C	0.75	3/2531 (0.1%)	0.74	0/3429
All	All	0.86	7/5065 (0.1%)	0.81	5/6862 (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	C	0	1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	225	GLU	CD-OE2	8.62	1.35	1.25
1	C	154	GLU	CD-OE2	7.78	1.34	1.25
1	C	225	GLU	CD-OE2	7.77	1.34	1.25
1	C	242	GLU	CD-OE2	7.74	1.34	1.25
1	A	242	GLU	CD-OE2	7.15	1.33	1.25
1	A	154	GLU	CD-OE2	5.66	1.31	1.25
1	A	129	CYS	CB-SG	-5.02	1.73	1.81

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	281	ARG	NE-CZ-NH1	7.74	124.17	120.30
1	A	281	ARG	NE-CZ-NH2	-6.83	116.89	120.30
1	A	69	VAL	CB-CA-C	-6.07	99.86	111.40
1	A	252	ARG	NE-CZ-NH2	-5.99	117.31	120.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	72	ASP	N-CA-CB	-5.12	101.38	110.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	94	GLU	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	2474	0	2363	68	2
1	C	2471	0	2359	73	1
2	A	1	0	0	1	0
3	A	104	0	0	14	0
3	C	29	0	0	9	0
All	All	5079	0	4722	141	2

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 15.

All (141) 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:166:ARG:HG3	1:A:166:ARG:HH11	1.10	1.09
1:C:17:ARG:HD3	3:C:342:HOH:O	1.52	1.06
1:C:219:GLN:HB2	1:C:314:PHE:CE2	1.99	0.98
1:A:94:GLU:HG2	1:A:96:SER:H	1.31	0.92
1:C:225:GLU:O	1:C:229:LEU:HD12	1.72	0.89
1:A:205:ARG:HD3	3:A:350:HOH:O	1.73	0.88
1:A:121:THR:O	1:A:281:ARG:NH2	2.09	0.85
1:A:34:LYS:HE3	1:A:34:LYS:HA	1.58	0.84
1:A:16:GLU:HB3	3:A:394:HOH:O	1.76	0.84
1:C:14:SER:O	1:C:18:LEU:HD13	1.77	0.82

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:31:LYS:HE3	1:C:72:ASP:OD2	1.79	0.82
1:A:166:ARG:HG3	1:A:166:ARG:NH1	1.89	0.79
1:A:166:ARG:HH11	1:A:166:ARG:CG	1.94	0.76
1:A:13:THR:HG23	3:A:420:HOH:O	1.84	0.76
1:C:52:TYR:O	1:C:56:THR:HG23	1.86	0.76
1:A:106:VAL:HG21	3:A:415:HOH:O	1.86	0.75
1:C:51:LYS:NZ	1:C:51:LYS:HB3	2.01	0.74
1:C:89:ASN:C	1:C:91:THR:H	1.92	0.71
1:C:124:LYS:HE2	3:C:324:HOH:O	1.90	0.71
1:C:161:LEU:HD23	1:C:169:GLY:HA2	1.72	0.70
1:A:13:THR:CG2	3:A:420:HOH:O	2.39	0.70
1:C:36:ILE:HD12	1:C:36:ILE:H	1.57	0.69
1:A:10:ASN:O	1:A:11:ASP:HB2	1.92	0.69
1:A:106:VAL:CG2	3:A:415:HOH:O	2.40	0.68
1:C:14:SER:HA	3:C:327:HOH:O	1.92	0.68
1:A:190:ARG:HD2	3:A:411:HOH:O	1.93	0.67
1:A:94:GLU:HG2	1:A:95:LEU:N	2.09	0.67
1:A:14:SER:HB3	1:A:17:ARG:HD3	1.75	0.67
1:C:220:VAL:HG22	1:C:313:SER:HB2	1.76	0.67
1:A:191:ASN:HB3	2:A:320:CL:CL	2.32	0.67
1:A:190:ARG:CD	3:A:411:HOH:O	2.42	0.66
1:A:242:GLU:HG3	1:A:263:LYS:O	1.96	0.66
1:C:146:LYS:HD3	1:C:152:LEU:HB2	1.77	0.65
1:C:89:ASN:O	1:C:91:THR:N	2.29	0.64
1:A:98:GLU:HG3	1:A:98:GLU:O	1.96	0.64
1:C:12:LEU:HA	3:C:342:HOH:O	1.98	0.64
1:C:271:VAL:HG23	1:C:279:LEU:HD23	1.79	0.64
1:C:118:GLY:O	1:C:146:LYS:HE3	2.00	0.61
1:C:296:ILE:HG23	1:C:306:VAL:HG22	1.83	0.60
1:C:89:ASN:HB3	3:C:334:HOH:O	2.02	0.59
1:A:297:LYS:NZ	3:A:370:HOH:O	2.35	0.58
1:A:251:TYR:CZ	1:A:294:ILE:HG13	2.38	0.58
1:A:36:ILE:O	1:A:40:ILE:HG13	2.04	0.58
1:C:215:ASP:OD2	1:C:319:ASN:HB2	2.04	0.57
1:A:94:GLU:HA	3:A:376:HOH:O	2.05	0.57
1:C:91:THR:HG23	3:C:334:HOH:O	2.03	0.57
1:C:296:ILE:CG2	1:C:306:VAL:HG22	2.36	0.56
1:A:35:ASN:ND2	1:A:38:GLU:H	2.05	0.55
1:A:159:GLU:OE2	1:A:188:HIS:ND1	2.38	0.55
1:A:86:ILE:O	1:A:87:ALA:C	2.44	0.55
1:A:101:LEU:O	1:A:102:ASN:HB2	2.08	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:83:THR:CG2	1:C:84:GLY:N	2.70	0.54
1:C:315:TYR:HB2	1:C:316:PRO:HD2	1.90	0.54
1:C:203:ARG:O	1:C:206:GLU:HG2	2.08	0.53
1:C:219:GLN:HB2	1:C:314:PHE:CD2	2.43	0.53
1:C:89:ASN:C	1:C:91:THR:N	2.61	0.53
1:A:161:LEU:HD22	1:A:169:GLY:HA2	1.91	0.52
1:C:83:THR:HG23	1:C:84:GLY:H	1.73	0.52
1:A:124:LYS:NZ	1:A:154:GLU:OE1	2.36	0.52
1:A:35:ASN:HD22	1:A:35:ASN:C	2.13	0.52
1:A:35:ASN:HD22	1:A:38:GLU:H	1.58	0.52
1:C:156:SER:HB2	1:C:193:TYR:HB3	1.92	0.52
1:A:59:LYS:O	1:A:60:ASN:HB3	2.09	0.52
1:A:106:VAL:HB	1:A:230:TYR:CD1	2.45	0.51
1:C:296:ILE:HG23	1:C:306:VAL:CG2	2.40	0.51
1:C:143:SER:O	1:C:147:ILE:HG12	2.10	0.51
1:A:190:ARG:HD3	3:A:411:HOH:O	2.08	0.51
1:C:79:LYS:O	1:C:83:THR:HB	2.10	0.51
1:C:136:SER:OG	1:C:238:SER:OG	2.26	0.51
1:A:83:THR:CG2	3:A:422:HOH:O	2.60	0.50
1:A:203:ARG:O	1:A:207:LYS:HG2	2.12	0.50
1:A:144:ILE:CG2	1:A:145:ILE:N	2.76	0.49
1:A:144:ILE:HG23	1:A:145:ILE:N	2.28	0.49
1:C:200:ARG:HB2	1:C:203:ARG:NH2	2.28	0.49
1:A:98:GLU:CG	1:A:98:GLU:O	2.61	0.49
1:A:241:LEU:HD23	1:A:309:LEU:HD13	1.94	0.49
1:A:69:VAL:HG22	1:A:287:GLY:O	2.13	0.48
1:C:308:GLY:O	1:C:310:TYR:N	2.45	0.48
1:C:139:SER:OG	1:C:281:ARG:NH1	2.47	0.48
1:A:277:TYR:C	1:A:277:TYR:CD2	2.87	0.47
1:A:142:GLU:OE2	1:A:154:GLU:HA	2.14	0.47
1:A:95:LEU:HD11	3:A:416:HOH:O	2.14	0.47
1:C:68:ASN:HB2	1:C:287:GLY:O	2.15	0.47
1:C:14:SER:CA	3:C:327:HOH:O	2.58	0.47
1:C:14:SER:C	3:C:327:HOH:O	2.52	0.47
1:C:146:LYS:HD2	1:C:150:GLY:C	2.35	0.47
1:C:106:VAL:O	1:C:108:ILE:HG23	2.14	0.47
1:A:242:GLU:HG3	1:A:263:LYS:C	2.35	0.46
1:A:106:VAL:HB	1:A:230:TYR:HD1	1.80	0.46
1:C:86:ILE:O	1:C:87:ALA:HB3	2.15	0.46
1:A:92:THR:HG22	3:A:377:HOH:O	2.15	0.46
1:C:251:TYR:OH	1:C:254:GLY:O	2.20	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:297:LYS:HG3	1:C:298:ARG:N	2.30	0.46
1:A:138:VAL:O	1:A:142:GLU:HG3	2.14	0.46
1:C:161:LEU:O	1:C:199:GLN:NE2	2.41	0.46
1:C:274:GLY:HA3	1:C:277:TYR:CE1	2.50	0.46
1:C:224:ASN:OD1	1:C:226:GLY:N	2.48	0.46
1:C:137:ALA:O	1:C:141:ILE:HG12	2.15	0.46
1:C:218:ARG:HG2	1:C:317:VAL:CG2	2.45	0.46
1:A:34:LYS:HB2	1:A:38:GLU:OE1	2.16	0.46
1:C:87:ALA:C	1:C:89:ASN:H	2.19	0.46
1:C:160:LEU:HD13	1:C:178:ALA:HB1	1.96	0.45
1:C:13:THR:O	1:C:13:THR:HG23	2.15	0.45
1:C:83:THR:HG23	1:C:84:GLY:N	2.32	0.45
1:A:108:ILE:HG13	1:A:273:TYR:OH	2.17	0.45
1:A:94:GLU:HG2	1:A:96:SER:N	2.14	0.45
1:C:144:ILE:O	1:C:148:ARG:HG2	2.16	0.45
1:A:101:LEU:HD11	1:A:220:VAL:HA	1.99	0.45
1:C:308:GLY:O	1:C:311:THR:N	2.50	0.44
1:C:184:GLN:HB3	1:C:185:TYR:CD2	2.52	0.44
1:C:238:SER:O	1:C:313:SER:HA	2.18	0.44
1:C:280:ILE:HB	1:C:294:ILE:HG23	1.98	0.44
1:A:68:ASN:HB2	1:A:287:GLY:O	2.17	0.44
1:A:72:ASP:HB3	1:A:73:MET:HG2	1.99	0.44
1:C:161:LEU:CD2	1:C:169:GLY:HA2	2.43	0.44
1:C:14:SER:O	1:C:18:LEU:CD1	2.58	0.43
1:A:115:ARG:HD3	1:A:293:TYR:CZ	2.53	0.43
1:C:142:GLU:HG2	1:C:153:ASN:O	2.18	0.43
1:C:174:TYR:HB2	1:C:177:SER:HB3	2.00	0.43
1:C:277:TYR:HA	1:C:296:ILE:O	2.19	0.43
1:A:59:LYS:O	1:A:60:ASN:CB	2.67	0.43
1:A:160:LEU:O	1:A:161:LEU:C	2.58	0.43
1:A:161:LEU:HD23	1:A:161:LEU:HA	1.80	0.42
1:C:126:GLN:HG3	1:C:127:GLY:O	2.20	0.42
1:C:28:LYS:HB3	1:C:28:LYS:HE2	1.82	0.42
1:A:45:ILE:HD13	1:A:45:ILE:HA	1.87	0.41
1:C:142:GLU:OE2	1:C:154:GLU:HA	2.19	0.41
1:C:203:ARG:HD2	3:C:348:HOH:O	2.20	0.41
1:A:179:LEU:HD23	1:A:179:LEU:HA	1.90	0.41
1:A:13:THR:OG1	1:A:17:ARG:NH1	2.44	0.41
1:C:155:TYR:HD2	1:C:188:HIS:O	2.04	0.41
1:A:34:LYS:CE	1:A:34:LYS:HA	2.37	0.41
1:A:69:VAL:HG22	1:A:287:GLY:C	2.40	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:135:PHE:O	1:C:139:SER:HB2	2.21	0.41
1:C:200:ARG:HB3	1:C:201:TYR:H	1.73	0.40
1:A:144:ILE:HD13	1:A:235:GLN:NE2	2.36	0.40
1:A:171:ASN:HA	1:A:171:ASN:HD22	1.71	0.40
1:C:51:LYS:HZ2	1:C:51:LYS:HB3	1.82	0.40
1:A:56:THR:C	1:A:58:LYS:H	2.25	0.40
1:A:11:ASP:O	1:A:18:LEU:HD13	2.21	0.40
1:C:309:LEU:HD12	1:C:309:LEU:HA	1.72	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:GLU:OE2	1:C:44:GLU:OE2[1_456]	2.13	0.07
1:A:11:ASP:OD1	1:A:98:GLU:OE2[2_545]	2.17	0.03

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	308/363 (85%)	280 (91%)	19 (6%)	9 (3%)	6	9
1	C	308/363 (85%)	261 (85%)	37 (12%)	10 (3%)	5	7
All	All	616/726 (85%)	541 (88%)	56 (9%)	19 (3%)	5	8

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	11	ASP
1	A	87	ALA
1	A	91	THR
1	A	92	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	98	GLU
1	A	102	ASN
1	A	105	ASP
1	C	55	GLU
1	C	90	TYR
1	C	96	SER
1	C	102	ASN
1	C	107	ASN
1	C	309	LEU
1	A	60	ASN
1	C	87	ALA
1	C	104	GLY
1	C	200	ARG
1	C	202	CYS
1	A	88	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	261/307 (85%)	224 (86%)	37 (14%)	4	6
1	C	260/307 (85%)	224 (86%)	36 (14%)	4	7
All	All	521/614 (85%)	448 (86%)	73 (14%)	4	7

All (73) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	ASP
1	A	12	LEU
1	A	17	ARG
1	A	18	LEU
1	A	27	LEU
1	A	32	ILE
1	A	34	LYS
1	A	35	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	50	LEU
1	A	54	ASP
1	A	67	LEU
1	A	69	VAL
1	A	72	ASP
1	A	81	LYS
1	A	83	THR
1	A	91	THR
1	A	92	THR
1	A	94	GLU
1	A	98	GLU
1	A	101	LEU
1	A	103	ASP
1	A	106	VAL
1	A	108	ILE
1	A	128	SER
1	A	161	LEU
1	A	166	ARG
1	A	180	GLN
1	A	193	TYR
1	A	205	ARG
1	A	217	VAL
1	A	221	GLN
1	A	241	LEU
1	A	246	LYS
1	A	257	VAL
1	A	303	SER
1	A	306	VAL
1	A	319	ASN
1	C	10	ASN
1	C	12	LEU
1	C	14	SER
1	C	15	THR
1	C	18	LEU
1	C	19	ILE
1	C	24	SER
1	C	32	ILE
1	C	36	ILE
1	C	51	LYS
1	C	54	ASP
1	C	55	GLU
1	C	56	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	67	LEU
1	C	69	VAL
1	C	83	THR
1	C	91	THR
1	C	93	THR
1	C	106	VAL
1	C	129	CYS
1	C	139	SER
1	C	143	SER
1	C	144	ILE
1	C	145	ILE
1	C	161	LEU
1	C	180	GLN
1	C	181	LEU
1	C	193	TYR
1	C	219	GLN
1	C	235	GLN
1	C	241	LEU
1	C	257	VAL
1	C	260	CYS
1	C	278	ILE
1	C	297	LYS
1	C	306	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	ASN
1	A	60	ASN
1	A	171	ASN
1	A	262	ASN
1	C	180	GLN
1	C	221	GLN
1	C	234	ASN
1	C	235	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is 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	310/363 (85%)	0.04	13 (4%) 40 33	30, 47, 111, 147	0
1	C	310/363 (85%)	0.23	14 (4%) 37 30	49, 74, 139, 182	0
All	All	620/726 (85%)	0.14	27 (4%) 38 31	30, 61, 128, 182	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	92	THR	6.8
1	C	106	VAL	6.3
1	A	102	ASN	6.2
1	A	101	LEU	5.8
1	C	104	GLY	5.2
1	A	103	ASP	5.1
1	C	105	ASP	3.6
1	A	92	THR	3.5
1	C	108	ILE	3.4
1	C	90	TYR	3.3
1	C	11	ASP	3.2
1	C	93	THR	3.2
1	A	105	ASP	3.1
1	C	91	THR	3.1
1	A	10	ASN	2.8
1	A	91	THR	2.8
1	C	96	SER	2.5
1	A	100	VAL	2.5
1	C	230	TYR	2.5
1	C	99	GLU	2.5
1	A	99	GLU	2.4
1	C	100	VAL	2.3
1	A	106	VAL	2.3
1	A	13	THR	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	104	GLY	2.2
1	C	37	ASP	2.1
1	A	12	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no 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
2	CL	A	320	1/1	0.90	0.23	4.26	97,97,97,97	0

6.5 Other polymers [i](#)

There are no such residues in this entry.