



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 28, 2016 – 07:46 AM EST

PDB ID : 5TXC
Title : AtxE2 Isopeptidase - APO
Authors : Chekan, J.R.; Nair, S.K.
Deposited on : 2016-11-16
Resolution : 2.40 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	:	unknown
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20028442
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)	:	rb-20028442

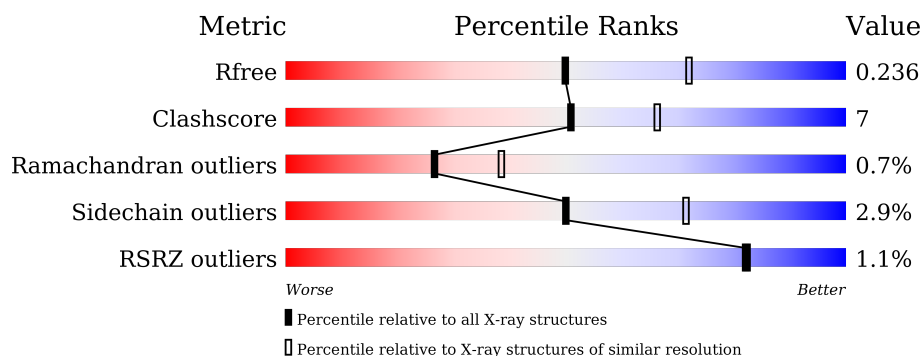
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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	705	
1	B	705	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10881 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 AtxE2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	653	Total	C	N	O	S	0	0	0
			5158	3245	936	962	15			
1	B	653	Total	C	N	O	S	0	0	0
			5158	3245	936	962	15			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	696	GLY	-	expression tag	UNP E8RUP5
A	697	SER	-	expression tag	UNP E8RUP5
A	698	ARG	-	expression tag	UNP E8RUP5
A	699	SER	-	expression tag	UNP E8RUP5
A	700	HIS	-	expression tag	UNP E8RUP5
A	701	HIS	-	expression tag	UNP E8RUP5
A	702	HIS	-	expression tag	UNP E8RUP5
A	703	HIS	-	expression tag	UNP E8RUP5
A	704	HIS	-	expression tag	UNP E8RUP5
A	705	HIS	-	expression tag	UNP E8RUP5
B	696	GLY	-	expression tag	UNP E8RUP5
B	697	SER	-	expression tag	UNP E8RUP5
B	698	ARG	-	expression tag	UNP E8RUP5
B	699	SER	-	expression tag	UNP E8RUP5
B	700	HIS	-	expression tag	UNP E8RUP5
B	701	HIS	-	expression tag	UNP E8RUP5
B	702	HIS	-	expression tag	UNP E8RUP5
B	703	HIS	-	expression tag	UNP E8RUP5
B	704	HIS	-	expression tag	UNP E8RUP5
B	705	HIS	-	expression tag	UNP E8RUP5

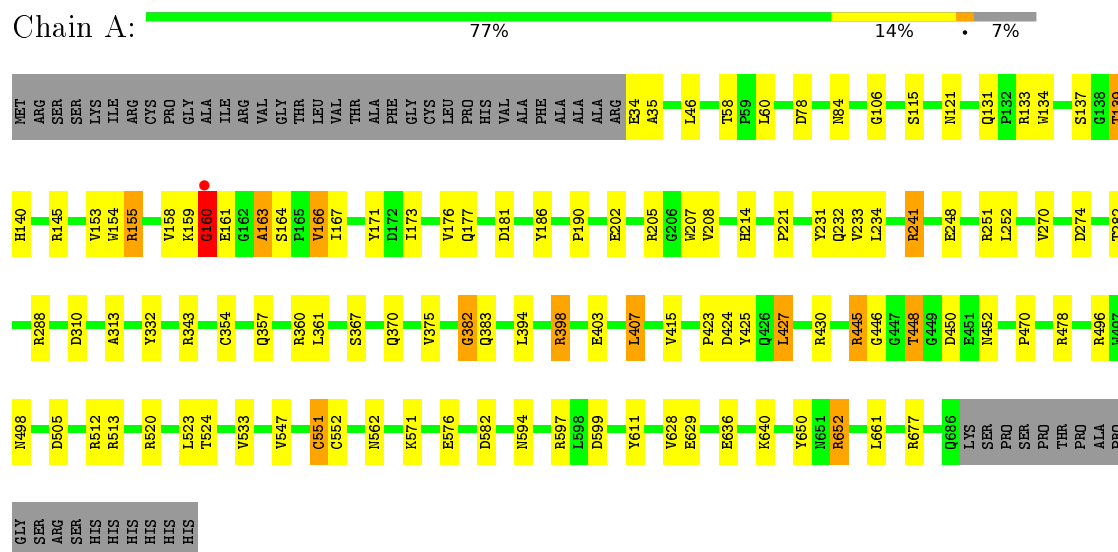
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	372	Total 372	O 372	0	0
2	B	193	Total 193	O 193	0	0

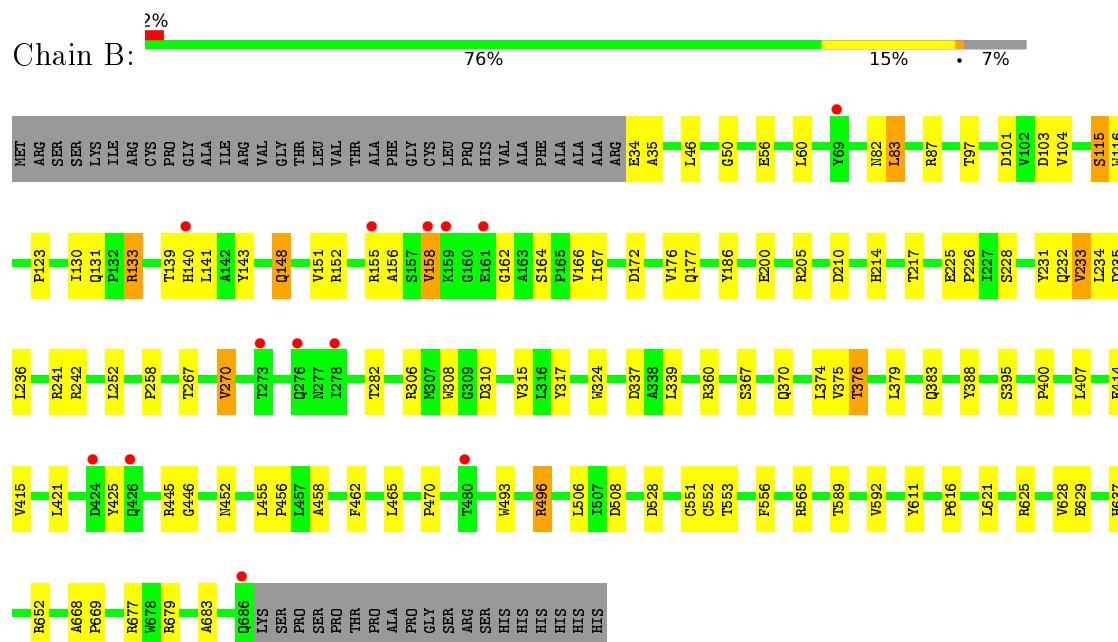
3 Residue-property plots

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: AtxE2



• Molecule 1: AtxE2



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	150.49Å 201.74Å 109.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.72 – 2.40 49.72 – 2.40	Depositor EDS
% Data completeness (in resolution range)	96.1 (49.72-2.40) 96.1 (49.72-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 2.39Å)	Xtriage
Refinement program	PHENIX (1.10 _2155: ???)	Depositor
R, R_{free}	0.194 , 0.236 0.192 , 0.236	Depositor DCC
R_{free} test set	3166 reflections (5.07%)	DCC
Wilson B-factor (Å ²)	39.0	Xtriage
Anisotropy	0.138	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 40.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10881	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.75% 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

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.46	1/5285 (0.0%)	0.71	7/7207 (0.1%)
1	B	0.40	0/5285	0.62	2/7207 (0.0%)
All	All	0.43	1/10570 (0.0%)	0.67	9/14414 (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	A	0	3
1	B	0	1
All	All	0	4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	58	THR	C-N	6.21	1.46	1.34

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	407	LEU	CA-CB-CG	7.10	131.63	115.30
1	A	445	ARG	NE-CZ-NH2	-6.44	117.08	120.30
1	A	398	ARG	NE-CZ-NH1	-6.16	117.22	120.30
1	B	235	ASP	CB-CG-OD2	5.40	123.16	118.30
1	A	78	ASP	CB-CG-OD1	5.24	123.02	118.30
1	A	448	THR	C-N-CA	-5.17	111.43	122.30
1	B	210	ASP	CB-CG-OD1	5.15	122.93	118.30
1	A	551	CYS	CA-C-N	-5.12	105.92	117.20
1	A	652	ARG	NE-CZ-NH1	-5.12	117.74	120.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	160	GLY	Peptide
1	A	382	GLY	Peptide
1	A	551	CYS	Peptide
1	B	551	CYS	Peptide

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5158	0	5061	66	0
1	B	5158	0	5061	78	0
2	A	372	0	0	19	1
2	B	193	0	0	19	0
All	All	10881	0	10122	141	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

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:205:ARG:O	1:B:652:ARG:NH2	2.09	0.85
1:A:205:ARG:NH1	2:A:803:HOH:O	2.10	0.84
1:A:652:ARG:NH1	2:A:805:HOH:O	2.13	0.81
1:A:448:THR:O	2:A:801:HOH:O	2.05	0.74
1:B:131:GLN:N	2:B:801:HOH:O	1.98	0.73
1:A:382:GLY:HA3	2:A:954:HOH:O	1.90	0.71
1:B:339:LEU:O	2:B:802:HOH:O	2.07	0.71
1:A:248:GLU:O	2:A:802:HOH:O	2.09	0.70
1:B:56:GLU:O	2:B:803:HOH:O	2.08	0.70
1:A:251:ARG:HB2	2:A:802:HOH:O	1.92	0.69
1:A:430:ARG:NH1	2:A:815:HOH:O	2.24	0.69
1:A:115:SER:HB3	1:A:121:ASN:ND2	2.08	0.68
1:B:465:LEU:HD21	1:B:506:LEU:HD13	1.76	0.68
1:B:34:GLU:O	2:B:804:HOH:O	2.13	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:679:ARG:NH2	2:B:811:HOH:O	2.27	0.67
1:B:270:VAL:HG13	1:B:282:THR:HB	1.76	0.66
1:B:228:SER:O	2:B:805:HOH:O	2.13	0.66
1:A:159:LYS:HA	1:A:161:GLU:OE1	1.96	0.66
1:A:498:ASN:OD1	2:A:807:HOH:O	2.14	0.66
1:B:130:ILE:HA	2:B:801:HOH:O	1.95	0.66
1:B:167:ILE:HG12	1:B:234:LEU:HD13	1.77	0.66
1:B:367:SER:HB3	1:B:370:GLN:HB3	1.77	0.66
1:B:677:ARG:NH2	2:B:815:HOH:O	2.29	0.65
1:A:599:ASP:OD1	2:A:804:HOH:O	2.13	0.65
1:B:167:ILE:HD13	1:B:186:TYR:CD2	2.35	0.62
1:B:133:ARG:HD2	1:B:176:VAL:O	2.00	0.62
1:B:103:ASP:OD2	1:B:143:TYR:OH	2.18	0.61
1:B:267:THR:OG1	2:B:806:HOH:O	2.16	0.60
1:A:357:GLN:NE2	2:A:825:HOH:O	2.34	0.59
1:A:652:ARG:NH2	1:B:205:ARG:O	2.36	0.59
1:B:306:ARG:HG2	1:B:308:TRP:CZ2	2.38	0.59
1:A:46:LEU:O	1:A:445:ARG:NH2	2.32	0.59
1:A:357:GLN:OE1	1:A:360:ARG:NH2	2.17	0.58
1:A:505:ASP:OD2	2:A:810:HOH:O	2.17	0.58
1:A:270:VAL:HG13	1:A:282:THR:HB	1.85	0.57
1:B:148:GLN:OE1	1:B:148:GLN:N	2.37	0.57
1:B:629:GLU:OE2	1:B:677:ARG:HD2	2.06	0.56
1:B:166:VAL:HG12	1:B:167:ILE:HG13	1.88	0.56
1:B:46:LEU:O	1:B:445:ARG:NH2	2.38	0.56
1:B:97:THR:N	2:B:821:HOH:O	2.36	0.55
1:A:403:GLU:OE2	1:A:513:ARG:NH2	2.40	0.54
1:A:163:ALA:HA	2:A:845:HOH:O	2.06	0.54
1:A:415:VAL:HG22	1:A:470:PRO:HB3	1.89	0.54
1:A:398:ARG:HG2	1:A:423:PRO:HB3	1.89	0.54
1:B:200:GLU:OE1	2:B:807:HOH:O	2.18	0.54
1:A:354:CYS:HB3	1:A:361:LEU:HG	1.90	0.54
1:A:398:ARG:NH2	1:A:424:ASP:OD1	2.41	0.54
1:A:367:SER:HB3	1:A:370:GLN:HB3	1.89	0.53
1:B:233:VAL:HG21	1:B:242:ARG:CZ	2.38	0.53
1:B:141:LEU:O	1:B:155:ARG:HA	2.08	0.53
1:A:450:ASP:OD1	1:A:650:TYR:OH	2.22	0.53
1:A:84:ASN:ND2	2:A:806:HOH:O	2.14	0.53
1:A:520:ARG:HD2	1:A:661:LEU:O	2.07	0.53
1:B:446:GLY:HA2	1:B:452:ASN:O	2.10	0.52
1:A:234:LEU:HD13	1:A:241:ARG:HD2	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:87:ARG:HD2	1:B:101:ASP:OD2	2.10	0.52
1:B:458:ALA:HA	1:B:462:PHE:O	2.10	0.52
1:B:317:TYR:CE1	1:B:379:LEU:HB3	2.45	0.51
1:B:177:GLN:HG2	1:B:252:LEU:HD12	1.92	0.51
1:B:186:TYR:CZ	1:B:232:GLN:HB2	2.45	0.51
1:B:233:VAL:HG13	1:B:242:ARG:HG3	1.91	0.51
1:A:523:LEU:HG	1:A:533:VAL:HG22	1.93	0.51
1:B:415:VAL:HG22	1:B:470:PRO:HB3	1.92	0.50
1:B:155:ARG:HH21	1:B:236:LEU:HD13	1.76	0.50
1:A:571:LYS:NZ	1:A:576:GLU:OE1	2.42	0.50
1:B:155:ARG:HH21	1:B:236:LEU:HD22	1.77	0.49
1:A:155:ARG:HG2	1:A:166:VAL:HG21	1.95	0.49
1:B:140:HIS:HA	1:B:156:ALA:O	2.12	0.49
1:B:324:TRP:CZ2	1:B:414:GLU:HG2	2.48	0.49
1:A:190:PRO:HD2	2:A:872:HOH:O	2.13	0.48
1:B:496:ARG:HH22	1:B:528:ASP:CG	2.16	0.48
1:A:231:TYR:CD2	1:A:252:LEU:HD23	2.49	0.48
1:B:508:ASP:OD1	2:B:808:HOH:O	2.19	0.47
1:B:337:ASP:HB2	2:B:898:HOH:O	2.13	0.47
1:B:103:ASP:OD1	1:B:104:VAL:N	2.47	0.47
1:A:231:TYR:HD2	1:A:252:LEU:HD23	1.79	0.47
1:A:445:ARG:HB3	2:A:896:HOH:O	2.15	0.47
1:A:153:VAL:HG11	1:A:176:VAL:HG21	1.97	0.47
1:B:233:VAL:HG21	1:B:242:ARG:NH2	2.30	0.47
1:A:139:THR:HG22	1:A:140:HIS:ND1	2.31	0.46
1:A:425:TYR:HE2	1:A:427:LEU:HD13	1.80	0.46
1:A:310:ASP:O	1:A:313:ALA:O	2.33	0.46
1:A:594:ASN:OD1	1:A:597:ARG:NH2	2.44	0.46
1:A:137:SER:OG	1:A:181:ASP:OD2	2.33	0.46
1:B:374:LEU:HB2	1:B:388:TYR:HB3	1.97	0.46
1:B:400:PRO:CG	1:B:455:LEU:HD22	2.46	0.46
1:B:115:SER:HB3	1:B:116:TRP:CD1	2.51	0.46
1:A:167:ILE:HD13	1:A:186:TYR:CD1	2.51	0.46
1:A:145:ARG:HB3	1:A:154:TRP:CD1	2.52	0.45
1:B:50:GLY:HA3	1:B:60:LEU:HD11	1.98	0.45
1:A:332:TYR:CE2	1:A:343:ARG:HB2	2.51	0.45
1:B:556:PHE:CD2	1:B:616:PRO:HG2	2.52	0.45
1:A:166:VAL:HG12	1:A:167:ILE:HG13	1.98	0.45
1:B:139:THR:O	1:B:158:VAL:HG22	2.17	0.45
1:B:565:ARG:NH2	2:B:812:HOH:O	2.26	0.45
1:B:156:ALA:HA	1:B:162:GLY:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:GLN:OE1	2:A:811:HOH:O	2.21	0.44
1:A:274:ASP:OD2	1:A:478:ARG:NH1	2.50	0.44
1:A:173:ILE:HD12	1:A:186:TYR:CD2	2.53	0.44
1:A:214:HIS:HD2	1:A:562:ASN:O	2.01	0.44
1:B:421:LEU:HD13	1:B:425:TYR:CD2	2.52	0.44
1:B:592:VAL:HG23	2:B:817:HOH:O	2.18	0.44
1:B:123:PRO:HG2	1:B:637:HIS:CE1	2.53	0.43
1:A:582:ASP:OD1	1:A:582:ASP:N	2.46	0.43
1:B:87:ARG:HB3	1:B:104:VAL:HG22	2.01	0.43
1:A:394:LEU:HA	1:A:394:LEU:HD12	1.69	0.43
1:B:231:TYR:CD2	1:B:252:LEU:HD23	2.54	0.43
1:B:258:PRO:HB2	1:B:306:ARG:HD2	2.01	0.43
1:A:106:GLY:HA3	2:A:987:HOH:O	2.18	0.43
1:B:621:LEU:HD13	1:B:628:VAL:HG11	2.01	0.43
1:B:360:ARG:HD2	1:B:376:THR:HG21	2.01	0.43
1:B:231:TYR:HD2	1:B:252:LEU:HD23	1.83	0.43
1:B:683:ALA:HA	2:B:850:HOH:O	2.18	0.43
1:A:202:GLU:HG2	2:A:803:HOH:O	2.19	0.42
1:B:152:ARG:HD2	2:B:882:HOH:O	2.19	0.42
1:A:134:TRP:CZ3	1:A:158:VAL:HG11	2.55	0.42
1:B:553:THR:HB	1:B:589:THR:HB	2.02	0.42
1:A:166:VAL:CG1	1:A:167:ILE:HG13	2.50	0.42
1:A:446:GLY:HA2	1:A:452:ASN:O	2.20	0.42
1:A:160:GLY:H	1:A:161:GLU:CD	2.23	0.42
1:B:225:GLU:OE1	1:B:226:PRO:HA	2.19	0.42
1:B:214:HIS:CE1	1:B:217:THR:HG23	2.54	0.41
1:B:400:PRO:HG3	1:B:455:LEU:HD22	2.02	0.41
1:B:625:ARG:NH1	2:B:834:HOH:O	2.53	0.41
1:B:82:ASN:O	1:B:83:LEU:O	2.38	0.41
1:B:155:ARG:HG2	1:B:166:VAL:CG2	2.49	0.41
2:A:855:HOH:O	1:B:652:ARG:NH1	2.53	0.41
1:A:629:GLU:OE2	1:A:677:ARG:HD2	2.21	0.41
1:B:151:VAL:HB	1:B:172:ASP:HB3	2.03	0.41
1:A:171:TYR:CD2	1:A:190:PRO:HB3	2.56	0.41
1:A:207:TRP:CE3	1:A:221:PRO:HD3	2.55	0.41
1:B:310:ASP:HB3	1:B:315:VAL:HG13	2.02	0.41
1:B:668:ALA:HA	1:B:669:PRO:HD3	1.99	0.41
1:A:636:GLU:OE1	1:A:640:LYS:HA	2.21	0.41
1:B:155:ARG:NH2	1:B:236:LEU:HD13	2.36	0.41
1:A:208:VAL:HG21	1:B:677:ARG:CZ	2.50	0.41
1:A:139:THR:HG22	1:A:140:HIS:CG	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:512:ARG:HH11	1:A:512:ARG:HD2	1.75	0.40
1:B:493:TRP:HA	2:B:843:HOH:O	2.20	0.40
1:A:524:THR:HA	1:A:547:VAL:O	2.21	0.40
1:B:455:LEU:HB2	1:B:456:PRO:HD3	2.02	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:837:HOH:O	2:A:1076:HOH:O[6_575]	1.92	0.28

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	651/705 (92%)	625 (96%)	21 (3%)	5 (1%)	24	35
1	B	651/705 (92%)	626 (96%)	21 (3%)	4 (1%)	30	43
All	All	1302/1410 (92%)	1251 (96%)	42 (3%)	9 (1%)	26	38

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	552	CYS
1	B	83	LEU
1	B	552	CYS
1	A	35	ALA
1	A	164	SER
1	B	35	ALA
1	B	164	SER
1	A	163	ALA
1	A	160	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	547/588 (93%)	529 (97%)	18 (3%)	45	66
1	B	547/588 (93%)	533 (97%)	14 (3%)	54	74
All	All	1094/1176 (93%)	1062 (97%)	32 (3%)	50	71

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	GLU
1	A	60	LEU
1	A	131	GLN
1	A	133	ARG
1	A	139	THR
1	A	155	ARG
1	A	166	VAL
1	A	232	GLN
1	A	233	VAL
1	A	241	ARG
1	A	288	ARG
1	A	375	VAL
1	A	383	GLN
1	A	407	LEU
1	A	427	LEU
1	A	496	ARG
1	A	611	TYR
1	A	628	VAL
1	B	115	SER
1	B	133	ARG
1	B	148	GLN
1	B	158	VAL
1	B	233	VAL
1	B	241	ARG
1	B	270	VAL
1	B	375	VAL
1	B	376	THR

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Mol	Chain	Res	Type
1	B	383	GLN
1	B	395	SER
1	B	407	LEU
1	B	496	ARG
1	B	611	TYR

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

Mol	Chain	Res	Type
1	A	73	GLN
1	A	121	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	653/705 (92%)	-0.40	1 (0%) 95 95	22, 33, 54, 86	0
1	B	653/705 (92%)	-0.05	13 (1%) 68 68	26, 46, 77, 109	0
All	All	1306/1410 (92%)	-0.23	14 (1%) 82 82	22, 39, 71, 109	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	160	GLY	4.0
1	B	158	VAL	3.9
1	B	273	THR	3.7
1	B	480	THR	3.4
1	B	276	GLN	3.0
1	B	426	GLN	2.7
1	B	424	ASP	2.7
1	B	159	LYS	2.5
1	B	278	ILE	2.4
1	B	69	TYR	2.4
1	B	161	GLU	2.3
1	B	140	HIS	2.3
1	B	155	ARG	2.2
1	B	686	GLN	2.2

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.