



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 2, 2017 – 05:53 PM EST

PDB ID : 5TUA
Title : structure of a Na⁺-selective mutant of two-pore channel from Arabidopsis thaliana AtTPC1
Authors : Guo, J.; Zeng, W.; Jiang, Y.
Deposited on : 2016-11-05
Resolution : 3.30 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

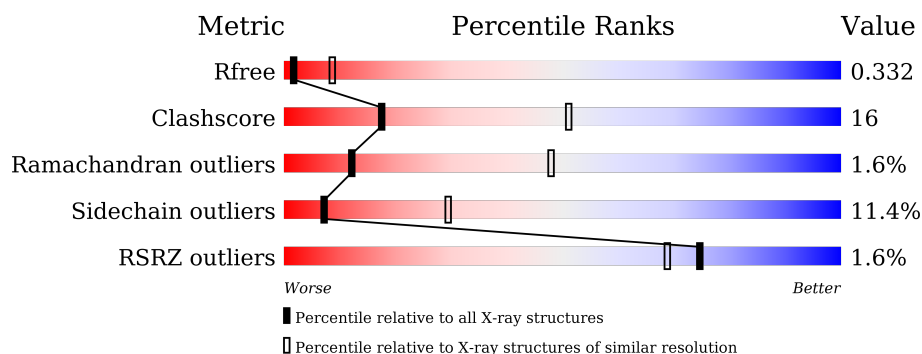
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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	2060 (3.40-3.20)
Clashscore	102246	1058 (3.38-3.22)
Ramachandran outliers	100387	1038 (3.38-3.22)
Sidechain outliers	100360	1037 (3.38-3.22)
RSRZ outliers	91569	2070 (3.40-3.20)

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	739	<div> <div></div> <div>47%</div> <div>33%</div> <div>15%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5126 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 Two pore calcium channel protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	627	Total	C	N	O	S	0	0	0
			5114	3399	787	907	21			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	265	ALA	SER	conflict	UNP Q94KI8
A	629	VAL	MET	conflict	UNP Q94KI8
A	630	ASN	GLY	conflict	UNP Q94KI8
A	734	SER	-	expression tag	UNP Q94KI8
A	735	THR	-	expression tag	UNP Q94KI8
A	736	ALA	-	expression tag	UNP Q94KI8
A	737	GLY	-	expression tag	UNP Q94KI8
A	738	LEU	-	expression tag	UNP Q94KI8
A	739	VAL	-	expression tag	UNP Q94KI8

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Ca	0	0
			2	2		

- Molecule 3 is BARIUM ION (three-letter code: BA) (formula: Ba).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	9	Total	Ba	0	0
			9	9		

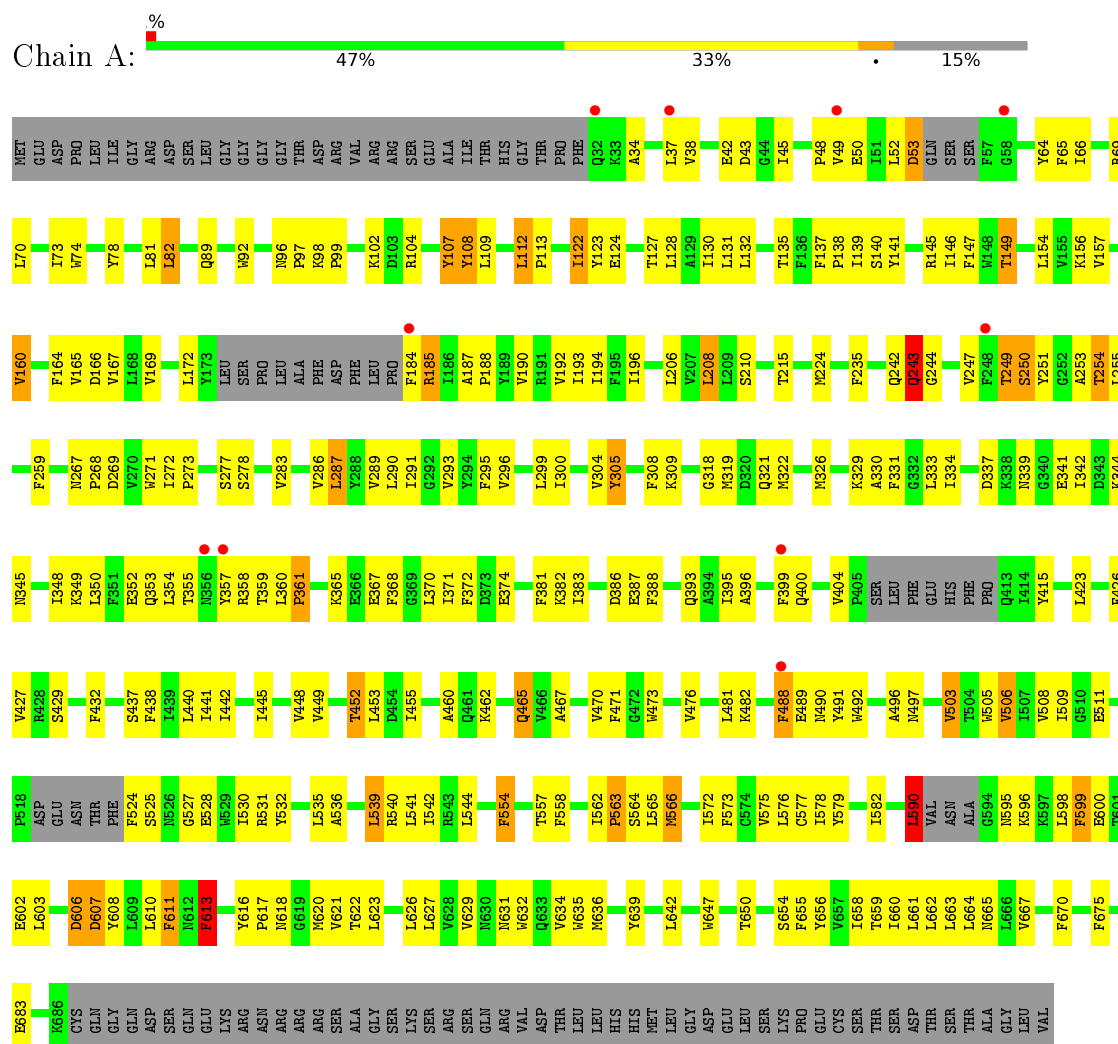
- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Na	0	0
			1	1		

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: Two pore calcium channel protein 1



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	88.53Å 156.37Å 217.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.59 – 3.30 44.59 – 3.29	Depositor EDS
% Data completeness (in resolution range)	62.8 (44.59-3.30) 62.8 (44.59-3.29)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.65 (at 3.32Å)	Xtriage
Refinement program	PHENIX (1.10_2155: ???)	Depositor
R, R_{free}	0.310 , 0.323 0.307 , 0.332	Depositor DCC
R_{free} test set	725 reflections (4.96%)	DCC
Wilson B-factor (Å ²)	25.1	Xtriage
Anisotropy	0.599	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 13.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.036 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.045 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.72	EDS
Total number of atoms	5126	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 6.14% 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: NA, CA, BA

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	0.32	0/5241	0.52	1/7128 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	590	LEU	CA-CB-CG	5.68	128.36	115.30

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	5114	0	5091	163	0
2	A	2	0	0	0	0
3	A	9	0	0	0	0
4	A	1	0	0	0	0
All	All	5126	0	5091	163	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 16.

All (163) 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:577:CYS:HA	1:A:620:MET:HE1	1.64	0.79
1:A:250:SER:OG	1:A:251:TYR:N	2.11	0.79
1:A:536:ALA:O	1:A:539:LEU:HB2	1.83	0.78
1:A:572:ILE:HD11	1:A:627:LEU:HD11	1.66	0.76
1:A:66:ILE:HD13	1:A:140:SER:HB2	1.68	0.74
1:A:603:LEU:HD13	1:A:606:ASP:HB2	1.68	0.73
1:A:613:PHE:H	1:A:618:ASN:HB3	1.53	0.73
1:A:367:GLU:HA	1:A:370:LEU:HB2	1.71	0.71
1:A:348:ILE:O	1:A:352:GLU:HB2	1.90	0.71
1:A:112:LEU:HD22	1:A:113:PRO:HD2	1.73	0.71
1:A:353:GLN:O	1:A:357:TYR:HB2	1.91	0.71
1:A:124:GLU:O	1:A:128:LEU:HB2	1.91	0.70
1:A:190:VAL:O	1:A:194:ILE:HB	1.92	0.70
1:A:595:ASN:O	1:A:599:PHE:HB2	1.92	0.69
1:A:145:ARG:O	1:A:149:THR:HB	1.94	0.68
1:A:38:VAL:HG21	1:A:353:GLN:HB2	1.74	0.67
1:A:367:GLU:O	1:A:371:ILE:HB	1.94	0.66
1:A:656:TYR:O	1:A:660:ILE:HB	1.95	0.66
1:A:370:LEU:O	1:A:374:GLU:HB3	1.95	0.66
1:A:344:LYS:NZ	1:A:345:ASN:OD1	2.29	0.65
1:A:74:TRP:O	1:A:78:TYR:CB	2.46	0.64
1:A:607:ASP:HA	1:A:610:LEU:HB3	1.81	0.62
1:A:49:VAL:HG22	1:A:322:MET:HE2	1.83	0.61
1:A:37:LEU:CD1	1:A:334:ILE:HB	2.31	0.60
1:A:259:PHE:HE2	1:A:632:TRP:HH2	1.47	0.60
1:A:354:LEU:HD11	1:A:395:ILE:HG21	1.82	0.60
1:A:235:PHE:HB2	1:A:254:THR:HG21	1.84	0.60
1:A:438:PHE:HA	1:A:441:ILE:HD12	1.84	0.60
1:A:81:LEU:HD11	1:A:127:THR:HA	1.83	0.59
1:A:157:VAL:HA	1:A:160:VAL:HB	1.83	0.59
1:A:350:LEU:HD21	1:A:388:PHE:CE1	2.36	0.59
1:A:53:ASP:OD1	1:A:53:ASP:N	2.35	0.59
1:A:249:THR:HB	1:A:253:ALA:HB2	1.84	0.59
1:A:334:ILE:HD11	1:A:349:LYS:HZ2	1.67	0.59
1:A:337:ASP:OD2	1:A:339:ASN:ND2	2.35	0.58
1:A:623:LEU:HD23	1:A:626:LEU:HD12	1.84	0.58
1:A:291:ILE:O	1:A:295:PHE:HB2	2.04	0.58
1:A:524:PHE:O	1:A:528:GLU:N	2.33	0.58
1:A:655:PHE:O	1:A:659:THR:HG22	2.03	0.58
1:A:470:VAL:HA	1:A:473:TRP:HD1	1.68	0.57
1:A:286:VAL:O	1:A:290:LEU:HB2	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:598:LEU:O	1:A:602:GLU:CB	2.53	0.57
1:A:342:ILE:O	1:A:382:LYS:HA	2.05	0.55
1:A:448:VAL:O	1:A:452:THR:HB	2.06	0.55
1:A:331:PHE:HB2	1:A:388:PHE:CD2	2.41	0.55
1:A:511:GLU:HA	1:A:531:ARG:NH1	2.21	0.55
1:A:404:VAL:HA	1:A:490:ASN:HD21	1.71	0.55
1:A:505:TRP:HA	1:A:508:VAL:HG22	1.89	0.54
1:A:563:PRO:O	1:A:566:MET:HB2	2.08	0.54
1:A:579:TYR:HA	1:A:582:ILE:HD12	1.88	0.54
1:A:488:PHE:HD1	1:A:489:GLU:HG2	1.72	0.54
1:A:43:ASP:O	1:A:48:PRO:HD2	2.08	0.54
1:A:122:ILE:HG13	1:A:123:TYR:N	2.21	0.54
1:A:606:ASP:N	1:A:606:ASP:OD1	2.35	0.54
1:A:370:LEU:O	1:A:374:GLU:CB	2.56	0.53
1:A:138:PRO:HA	1:A:141:TYR:CE2	2.43	0.53
1:A:139:ILE:HG12	1:A:147:PHE:CG	2.43	0.53
1:A:107:TYR:OH	1:A:596:LYS:HG2	2.09	0.53
1:A:300:ILE:O	1:A:304:VAL:HG23	2.08	0.53
1:A:426:PHE:O	1:A:429:SER:OG	2.27	0.53
1:A:449:VAL:O	1:A:452:THR:HG22	2.08	0.53
1:A:109:LEU:HD22	1:A:611:PHE:HB2	1.91	0.52
1:A:45:ILE:HG12	1:A:396:ALA:HB2	1.92	0.52
1:A:208:LEU:HD21	1:A:308:PHE:HD1	1.74	0.52
1:A:34:ALA:HB3	1:A:353:GLN:HE22	1.75	0.52
1:A:165:VAL:O	1:A:169:VAL:HG23	2.10	0.52
1:A:112:LEU:HG	1:A:617:PRO:HG2	1.92	0.52
1:A:272:ILE:HG22	1:A:273:PRO:HD3	1.90	0.51
1:A:329:LYS:O	1:A:333:LEU:HG	2.10	0.51
1:A:554:PHE:HD1	1:A:554:PHE:N	2.07	0.51
1:A:539:LEU:HD23	1:A:542:ILE:HD13	1.92	0.51
1:A:138:PRO:HG2	1:A:147:PHE:HE2	1.75	0.51
1:A:267:ASN:O	1:A:271:TRP:NE1	2.43	0.51
1:A:166:ASP:OD2	1:A:187:ALA:HB2	2.11	0.50
1:A:437:SER:O	1:A:440:LEU:HG	2.11	0.50
1:A:554:PHE:CD1	1:A:554:PHE:N	2.76	0.50
1:A:393:GLN:HA	1:A:396:ALA:HB3	1.92	0.50
1:A:131:LEU:O	1:A:135:THR:HG23	2.11	0.50
1:A:453:LEU:HD13	1:A:460:ALA:HB3	1.94	0.50
1:A:164:PHE:O	1:A:167:VAL:HG12	2.12	0.50
1:A:330:ALA:O	1:A:333:LEU:N	2.43	0.49
1:A:467:ALA:O	1:A:470:VAL:HG12	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:664:LEU:HA	1:A:667:VAL:HG22	1.94	0.49
1:A:452:THR:HA	1:A:455:ILE:HG12	1.94	0.49
1:A:89:GLN:HG3	1:A:188:PRO:HB3	1.95	0.49
1:A:42:GLU:HB2	1:A:358:ARG:HG3	1.93	0.49
1:A:344:LYS:HB2	1:A:381:PHE:CB	2.43	0.49
1:A:564:SER:HB2	1:A:670:PHE:HZ	1.77	0.49
1:A:558:PHE:O	1:A:562:ILE:HG13	2.12	0.49
1:A:359:THR:HG21	1:A:400:GLN:HA	1.94	0.49
1:A:70:LEU:HB2	1:A:137:PHE:CE1	2.48	0.48
1:A:395:ILE:O	1:A:400:GLN:HB2	2.13	0.48
1:A:527:GLY:HA2	1:A:530:ILE:HG22	1.95	0.48
1:A:208:LEU:HD21	1:A:308:PHE:CD1	2.48	0.48
1:A:573:PHE:O	1:A:576:LEU:HB2	2.14	0.48
1:A:596:LYS:O	1:A:600:GLU:HG2	2.14	0.48
1:A:81:LEU:HD13	1:A:130:ILE:HD12	1.96	0.48
1:A:503:VAL:O	1:A:506:VAL:HG12	2.14	0.48
1:A:438:PHE:O	1:A:442:ILE:HG22	2.14	0.48
1:A:602:GLU:O	1:A:603:LEU:HD23	2.13	0.48
1:A:622:THR:OG1	1:A:635:TRP:NE1	2.47	0.47
1:A:250:SER:O	1:A:254:THR:HB	2.14	0.47
1:A:470:VAL:HA	1:A:473:TRP:CD1	2.47	0.47
1:A:590:LEU:HD22	1:A:642:LEU:HD22	1.95	0.47
1:A:337:ASP:N	1:A:337:ASP:OD1	2.35	0.47
1:A:187:ALA:HB3	1:A:188:PRO:HD3	1.97	0.47
1:A:135:THR:HG22	1:A:156:LYS:HE2	1.97	0.47
1:A:65:PHE:O	1:A:69:ARG:HG3	2.15	0.46
1:A:616:TYR:CE2	1:A:620:MET:HE2	2.51	0.46
1:A:37:LEU:HG	1:A:330:ALA:HB1	1.97	0.46
1:A:268:PRO:HG2	1:A:269:ASP:OD1	2.15	0.46
1:A:383:ILE:HD11	1:A:387:GLU:HG3	1.97	0.46
1:A:462:LYS:HA	1:A:465:GLN:HB3	1.96	0.46
1:A:496:ALA:HB1	1:A:540:ARG:HD2	1.98	0.46
1:A:243:GLN:NE2	1:A:271:TRP:HA	2.30	0.46
1:A:289:VAL:O	1:A:293:VAL:HB	2.15	0.46
1:A:506:VAL:HA	1:A:509:ILE:HG22	1.98	0.46
1:A:576:LEU:HD22	1:A:623:LEU:HB3	1.97	0.46
1:A:45:ILE:HG23	1:A:396:ALA:HB1	1.98	0.45
1:A:481:LEU:HD12	1:A:482:LYS:N	2.30	0.45
1:A:613:PHE:HB3	1:A:639:TYR:OH	2.17	0.45
1:A:334:ILE:HD12	1:A:350:LEU:HD22	1.99	0.44
1:A:650:THR:O	1:A:654:SER:HB3	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112:LEU:HD22	1:A:113:PRO:CD	2.46	0.44
1:A:355:THR:HA	1:A:361:PRO:HD2	1.99	0.44
1:A:78:TYR:O	1:A:82:LEU:HG	2.16	0.44
1:A:631:ASN:O	1:A:634:VAL:HG22	2.18	0.44
1:A:650:THR:O	1:A:654:SER:CB	2.66	0.44
1:A:575:VAL:HG21	1:A:663:LEU:HD11	1.98	0.44
1:A:247:VAL:HG13	1:A:254:THR:HB	2.00	0.43
1:A:66:ILE:O	1:A:69:ARG:HB2	2.18	0.43
1:A:146:ILE:O	1:A:149:THR:HG22	2.18	0.43
1:A:96:ASN:CG	1:A:97:PRO:HD3	2.38	0.43
1:A:554:PHE:HD1	1:A:554:PHE:H	1.67	0.43
1:A:296:VAL:O	1:A:300:ILE:HG13	2.18	0.43
1:A:661:LEU:O	1:A:665:ASN:ND2	2.36	0.43
1:A:361:PRO:HA	1:A:399:PHE:CD1	2.53	0.42
1:A:473:TRP:O	1:A:476:VAL:HG12	2.18	0.42
1:A:70:LEU:O	1:A:74:TRP:HB3	2.19	0.42
1:A:445:ILE:O	1:A:448:VAL:HG22	2.19	0.42
1:A:109:LEU:HD22	1:A:611:PHE:HD2	1.84	0.42
1:A:193:ILE:O	1:A:196:ILE:HG12	2.18	0.42
1:A:138:PRO:HG2	1:A:147:PHE:CE2	2.54	0.42
1:A:365:LYS:H	1:A:368:PHE:HB2	1.85	0.42
1:A:305:TYR:HE1	1:A:309:LYS:HE2	1.84	0.42
1:A:283:VAL:O	1:A:287:LEU:HB2	2.19	0.42
1:A:52:LEU:HD23	1:A:52:LEU:HA	1.91	0.42
1:A:491:TYR:CE1	1:A:497:ASN:HB3	2.55	0.42
1:A:453:LEU:HD23	1:A:453:LEU:HA	1.92	0.42
1:A:38:VAL:HG11	1:A:353:GLN:O	2.19	0.41
1:A:37:LEU:HD13	1:A:334:ILE:HB	1.99	0.41
1:A:557:THR:HG21	1:A:675:PHE:HB2	2.02	0.41
1:A:530:ILE:HG23	1:A:531:ARG:HG2	2.03	0.41
1:A:242:GLN:O	1:A:244:GLY:N	2.53	0.41
1:A:341:GLU:HA	1:A:383:ILE:O	2.20	0.41
1:A:318:GLY:O	1:A:321:GLN:HB2	2.21	0.41
1:A:130:ILE:H	1:A:130:ILE:HG13	1.68	0.41
1:A:432:PHE:CZ	1:A:481:LEU:HD11	2.56	0.41
1:A:184:PHE:CD1	1:A:185:ARG:HG2	2.56	0.40
1:A:172:LEU:HA	1:A:172:LEU:HD23	1.94	0.40
1:A:541:LEU:O	1:A:544:LEU:HB3	2.20	0.40
1:A:658:ILE:O	1:A:662:LEU:HB3	2.22	0.40
1:A:192:VAL:HG21	1:A:578:ILE:HG13	2.03	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	615/739 (83%)	561 (91%)	44 (7%)	10 (2%)	12	48

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	108	TYR
1	A	361	PRO
1	A	629	VAL
1	A	249	THR
1	A	613	PHE
1	A	104	ARG
1	A	243	GLN
1	A	563	PRO
1	A	647	TRP
1	A	99	PRO

5.3.2 Protein sidechains [i](#)

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	551/660 (84%)	488 (89%)	63 (11%)	7	29

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	GLU

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Mol	Chain	Res	Type
1	A	53	ASP
1	A	64	TYR
1	A	73	ILE
1	A	82	LEU
1	A	92	TRP
1	A	98	LYS
1	A	102	LYS
1	A	107	TYR
1	A	108	TYR
1	A	112	LEU
1	A	122	ILE
1	A	132	LEU
1	A	149	THR
1	A	154	LEU
1	A	160	VAL
1	A	185	ARG
1	A	206	LEU
1	A	208	LEU
1	A	210	SER
1	A	215	THR
1	A	224	MET
1	A	243	GLN
1	A	250	SER
1	A	254	THR
1	A	255	LEU
1	A	277	SER
1	A	278	SER
1	A	287	LEU
1	A	299	LEU
1	A	305	TYR
1	A	319	MET
1	A	326	MET
1	A	360	LEU
1	A	372	PHE
1	A	386	ASP
1	A	415	TYR
1	A	423	LEU
1	A	427	VAL
1	A	452	THR
1	A	465	GLN
1	A	471	PHE
1	A	488	PHE

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Mol	Chain	Res	Type
1	A	492	TRP
1	A	503	VAL
1	A	506	VAL
1	A	525	SER
1	A	532	TYR
1	A	535	LEU
1	A	539	LEU
1	A	554	PHE
1	A	565	LEU
1	A	566	MET
1	A	590	LEU
1	A	599	PHE
1	A	606	ASP
1	A	607	ASP
1	A	608	TYR
1	A	611	PHE
1	A	613	PHE
1	A	621	VAL
1	A	636	MET
1	A	683	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	267	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 12 ligands modelled in this entry, 12 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 [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	627/739 (84%)	-0.14	10 (1%) 74 69	16, 67, 114, 162	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	37	LEU	3.3
1	A	32	GLN	2.9
1	A	184	PHE	2.5
1	A	357	TYR	2.5
1	A	248	PHE	2.3
1	A	399	PHE	2.3
1	A	49	VAL	2.1
1	A	356	ASN	2.1
1	A	58	GLY	2.1
1	A	488	PHE	2.1

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	CA	A	801	1/1	0.92	0.05	-2.97	107,107,107,107	0
2	CA	A	802	1/1	0.91	0.06	-3.29	79,79,79,79	0
3	BA	A	808	1/1	0.96	0.05	-	136,136,136,136	0
3	BA	A	806	1/1	0.99	0.08	-	108,108,108,108	0
3	BA	A	807	1/1	0.89	0.14	-	156,156,156,156	0
3	BA	A	805	1/1	0.95	0.36	-	124,124,124,124	1
3	BA	A	810	1/1	0.97	0.34	-	118,118,118,118	1
3	BA	A	804	1/1	0.98	0.09	-	87,87,87,87	0
3	BA	A	809	1/1	0.88	0.20	-	180,180,180,180	0
3	BA	A	811	1/1	0.90	0.31	-	168,168,168,168	1
4	NA	A	812	1/1	0.93	0.11	-	27,27,27,27	1
3	BA	A	803	1/1	0.98	0.15	-	57,57,57,57	0

6.5 Other polymers [i](#)

There are no such residues in this entry.