Full wwPDB X-ray Structure Validation Report ⓘ

Feb 25, 2024 – 05:27 AM EST

PDB ID	:	5E1J
Title	:	Structure of voltage-gated two-pore channel TPC1 from Arabidopsis thaliana
Authors	:	Guo, J.; Zeng, W.; Chen, Q.; Lee, C.; Chen, L.; Yang, Y.; Jiang, Y.
Deposited on	:	2015-09-29
Resolution	:	3.31 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 types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references ⓘ) were used in the production of this report:

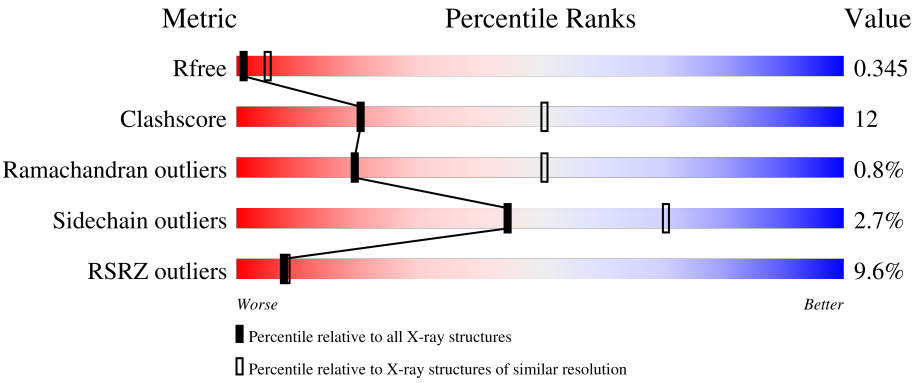
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.31 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	1149 (3.34-3.26)
Clashscore	141614	1205 (3.34-3.26)
Ramachandran outliers	138981	1183 (3.34-3.26)
Sidechain outliers	138945	1182 (3.34-3.26)
RSRZ outliers	127900	1115 (3.34-3.26)

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 <=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	741	8%, 62%, 20%, ·, 17%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 4964 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	616	Total	C	N	O	S	0	0	0
			4949	3286	762	880	21			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
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
A	740	PRO	-	expression tag	UNP Q94KI8
A	741	ARG	-	expression tag	UNP Q94KI8

- Molecule 2 is BARIUM ION (three-letter code: BA) (formula: Ba).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	8	Total	Ba	0	1
			9	9		

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 4 is water.

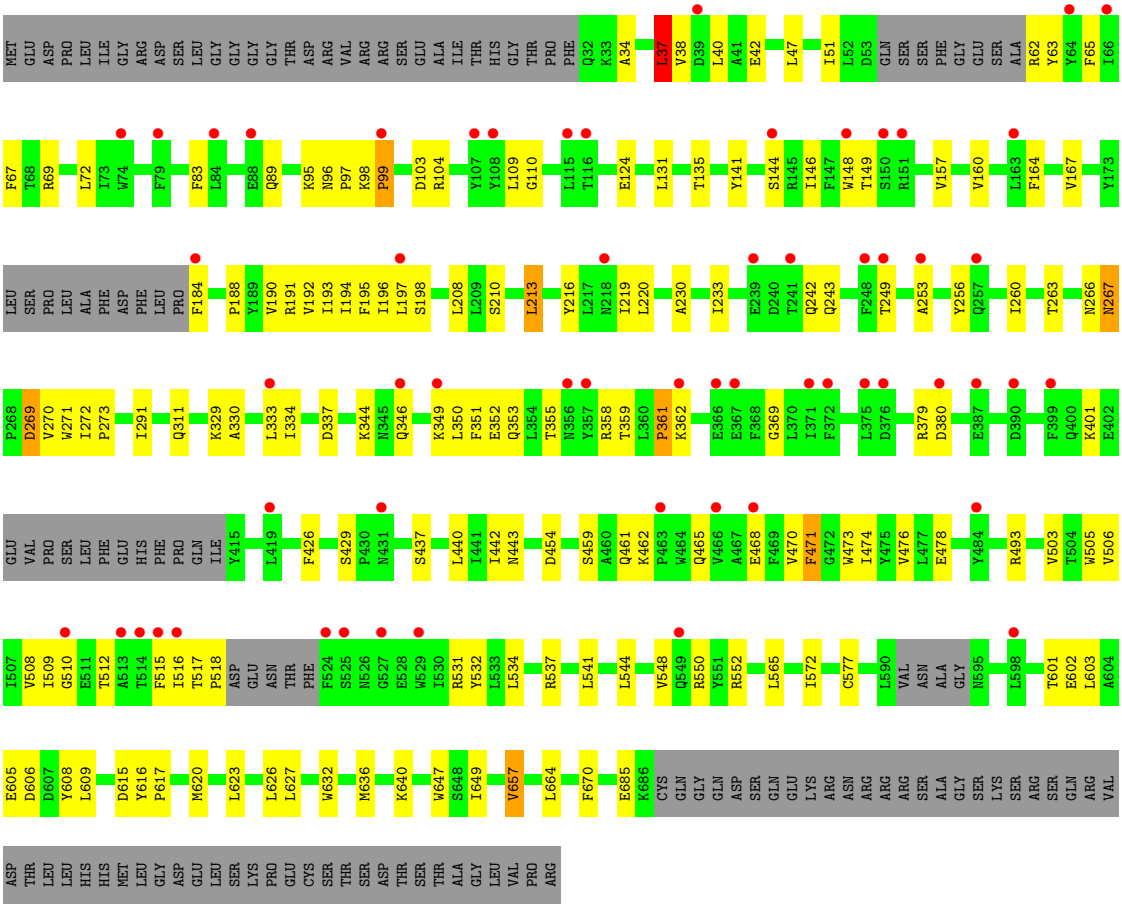
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	4	Total	O	0	0
			4	4		

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: Two pore calcium channel protein 1

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	88.43Å 158.85Å 217.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.71 – 3.31 39.71 – 3.31	Depositor EDS
% Data completeness (in resolution range)	77.3 (39.71-3.31) 77.6 (39.71-3.31)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$< I/\sigma(I) >$ [1]	19.97 (at 3.32Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.325 , 0.332 0.333 , 0.345	Depositor DCC
R_{free} test set	914 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å^2)	75.4	Xtriage
Anisotropy	0.791	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.27 , 30.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.022 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.029 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o,F_c correlation	0.83	EDS
Total number of atoms	4964	wwPDB-VP
Average B, all atoms (Å^2)	91.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.45% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\| > 5$	RMSZ	$\#\|Z\| > 5$
1	A	0.25	0/5067	0.44	2/6894 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	37	LEU	CB-CG-CD1	6.56	122.15	111.00
1	A	267	ASN	C-N-CD	5.08	139.08	128.40

There are no chirality outliers.

There are no planarity outliers.

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	4949	0	4875	122	0
2	A	9	0	0	0	0
3	A	2	0	0	0	0
4	A	4	0	0	0	0
All	All	4964	0	4875	122	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 12.

All (122) 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:63:TYR:O	1:A:67:PHE:HD2	1.06	1.29
1:A:63:TYR:O	1:A:67:PHE:CD2	1.85	1.28
1:A:272:ILE:CG2	1:A:273:PRO:HD3	1.74	1.16
1:A:146:ILE:O	1:A:149:THR:HG22	1.49	1.10
1:A:272:ILE:HG23	1:A:273:PRO:HD3	1.38	1.03
1:A:266:ASN:ND2	1:A:270:VAL:HB	1.82	0.94
1:A:272:ILE:HG22	1:A:273:PRO:HD3	1.47	0.93
1:A:266:ASN:HD22	1:A:270:VAL:HB	1.33	0.91
1:A:63:TYR:HB3	1:A:67:PHE:HE2	1.35	0.89
1:A:37:LEU:HD13	1:A:330:ALA:HB1	1.54	0.89
1:A:272:ILE:HG23	1:A:273:PRO:CD	2.06	0.84
1:A:63:TYR:HB3	1:A:67:PHE:CE2	2.11	0.84
1:A:63:TYR:C	1:A:67:PHE:HD2	1.81	0.83
1:A:269:ASP:O	1:A:272:ILE:HG22	1.79	0.83
1:A:131:LEU:HD21	1:A:194:ILE:CD1	2.12	0.80
1:A:272:ILE:CG2	1:A:273:PRO:CD	2.60	0.78
1:A:195:PHE:O	1:A:198:SER:OG	2.03	0.76
1:A:37:LEU:HD23	1:A:334:ILE:HB	1.72	0.71
1:A:63:TYR:CB	1:A:67:PHE:HE2	2.02	0.71
1:A:459:SER:HB2	1:A:461:GLN:HB2	1.71	0.71
1:A:63:TYR:C	1:A:67:PHE:CD2	2.60	0.71
1:A:213:LEU:HA	1:A:216:TYR:HB3	1.73	0.70
1:A:38:VAL:HG21	1:A:353:GLN:HB3	1.73	0.69
1:A:603:LEU:H	1:A:606:ASP:HB3	1.60	0.66
1:A:131:LEU:HD21	1:A:194:ILE:HD11	1.77	0.66
1:A:190:VAL:O	1:A:194:ILE:HG22	1.96	0.66
1:A:512:THR:HG23	1:A:516:ILE:HD12	1.78	0.65
1:A:197:LEU:C	1:A:197:LEU:HD12	2.17	0.64
1:A:266:ASN:ND2	1:A:270:VAL:CB	2.61	0.63
1:A:267:ASN:O	1:A:271:TRP:NE1	2.32	0.62
1:A:69:ARG:HA	1:A:72:LEU:HD23	1.82	0.62
1:A:193:ILE:O	1:A:196:ILE:HG12	2.00	0.62
1:A:503:VAL:HG13	1:A:534:LEU:HD12	1.83	0.61
1:A:548:VAL:HG12	1:A:550:ARG:H	1.66	0.60
1:A:216:TYR:OH	1:A:657:VAL:O	2.19	0.60
1:A:191:ARG:O	1:A:194:ILE:CG2	2.50	0.60
1:A:601:THR:HG22	1:A:602:GLU:HG3	1.83	0.59
1:A:606:ASP:HA	1:A:609:LEU:HB2	1.85	0.59
1:A:623:LEU:HD23	1:A:626:LEU:HD12	1.84	0.59
1:A:249:THR:HB	1:A:253:ALA:HB2	1.84	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:334:ILE:HD11	1:A:349:LYS:HZ2	1.69	0.58
1:A:552:ARG:NH1	1:A:685:GLU:OE2	2.36	0.57
1:A:191:ARG:O	1:A:194:ILE:HG22	2.03	0.57
1:A:135:THR:HG21	1:A:160:VAL:HG21	1.84	0.57
1:A:164:PHE:HA	1:A:167:VAL:HG12	1.89	0.55
1:A:197:LEU:HD12	1:A:198:SER:N	2.22	0.55
1:A:351:PHE:O	1:A:355:THR:HG23	2.08	0.54
1:A:465:GLN:O	1:A:468:GLU:HG2	2.08	0.53
1:A:95:LYS:NZ	1:A:184:PHE:O	2.32	0.52
1:A:454:ASP:OD2	1:A:461:GLN:NE2	2.43	0.52
1:A:572:ILE:HD11	1:A:627:LEU:HD11	1.92	0.52
1:A:192:VAL:O	1:A:196:ILE:HG23	2.09	0.51
1:A:42:GLU:HG2	1:A:358:ARG:HD2	1.92	0.50
1:A:334:ILE:HD11	1:A:349:LYS:NZ	2.26	0.50
1:A:505:TRP:HA	1:A:508:VAL:HG22	1.94	0.50
1:A:470:VAL:HA	1:A:473:TRP:HD1	1.77	0.49
1:A:37:LEU:HG	1:A:350:LEU:CD1	2.41	0.49
1:A:640:LYS:HZ2	1:A:640:LYS:HB3	1.77	0.49
1:A:208:LEU:HA	1:A:311:GLN:HG2	1.94	0.49
1:A:210:SER:HA	1:A:213:LEU:HD23	1.95	0.48
1:A:37:LEU:HG	1:A:350:LEU:HD11	1.95	0.48
1:A:577:CYS:HA	1:A:620:MET:HE1	1.96	0.48
1:A:89:GLN:HG3	1:A:188:PRO:HB3	1.96	0.48
1:A:426:PHE:O	1:A:429:SER:OG	2.29	0.48
1:A:565:LEU:HD13	1:A:670:PHE:HD2	1.79	0.48
1:A:344:LYS:HE3	1:A:369:GLY:HA2	1.96	0.47
1:A:38:VAL:HG21	1:A:353:GLN:CB	2.44	0.47
1:A:110:GLY:HA2	1:A:615:ASP:HB2	1.97	0.47
1:A:478:GLU:OE1	1:A:537:ARG:NH1	2.47	0.47
1:A:103:ASP:HA	1:A:104:ARG:HA	1.63	0.47
1:A:242:GLN:HA	1:A:243:GLN:HA	1.62	0.46
1:A:602:GLU:HA	1:A:603:LEU:HA	1.62	0.46
1:A:329:LYS:HE3	1:A:333:LEU:HD11	1.98	0.46
1:A:462:LYS:HA	1:A:465:GLN:HG2	1.98	0.46
1:A:355:THR:HG22	1:A:361:PRO:HG2	1.98	0.46
1:A:37:LEU:O	1:A:40:LEU:N	2.49	0.45
1:A:230:ALA:O	1:A:233:ILE:HG22	2.16	0.45
1:A:337:ASP:OD1	1:A:337:ASP:N	2.45	0.45
1:A:63:TYR:CA	1:A:67:PHE:CE2	2.99	0.45
1:A:98:LYS:HB3	1:A:99:PRO:HD2	1.99	0.45
1:A:34:ALA:O	1:A:38:VAL:HG23	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:359:THR:HG22	1:A:401:LYS:HA	1.98	0.45
1:A:379:ARG:CB	1:A:380:ASP:HB3	2.47	0.44
1:A:510:GLY:C	1:A:531:ARG:HH12	2.20	0.44
1:A:196:ILE:HG13	1:A:197:LEU:N	2.32	0.44
1:A:640:LYS:HB3	1:A:640:LYS:NZ	2.31	0.44
1:A:361:PRO:HB2	1:A:362:LYS:H	1.64	0.44
1:A:63:TYR:CA	1:A:67:PHE:HE2	2.29	0.44
1:A:197:LEU:C	1:A:197:LEU:CD1	2.85	0.44
1:A:605:GLU:O	1:A:609:LEU:N	2.45	0.44
1:A:62:ARG:HG2	1:A:65:PHE:HB2	2.00	0.43
1:A:191:ARG:O	1:A:194:ILE:HG23	2.17	0.43
1:A:352:GLU:O	1:A:355:THR:OG1	2.30	0.43
1:A:256:TYR:O	1:A:260:ILE:HG13	2.18	0.43
1:A:96:ASN:HB3	1:A:97:PRO:HD3	1.99	0.43
1:A:616:TYR:HB3	1:A:617:PRO:HD3	2.01	0.43
1:A:468:GLU:HA	1:A:471:PHE:HB2	1.99	0.43
1:A:148:TRP:HZ3	1:A:157:VAL:HG11	1.83	0.43
1:A:473:TRP:HA	1:A:476:VAL:HG12	2.01	0.43
1:A:263:THR:HG21	1:A:632:TRP:HE1	1.83	0.43
1:A:517:THR:HB	1:A:518:PRO:HD3	2.01	0.42
1:A:269:ASP:N	1:A:269:ASP:OD1	2.50	0.42
1:A:493:ARG:HA	1:A:493:ARG:HD3	1.85	0.42
1:A:440:LEU:HD11	1:A:541:LEU:HD13	2.01	0.42
1:A:506:VAL:HA	1:A:509:ILE:HG22	2.01	0.42
1:A:636:MET:HE2	1:A:649:ILE:HB	2.02	0.42
1:A:63:TYR:CB	1:A:67:PHE:CE2	2.88	0.41
1:A:67:PHE:HB2	1:A:141:TYR:HB3	2.02	0.41
1:A:144:SER:O	1:A:148:TRP:HD1	2.04	0.41
1:A:220:LEU:HD23	1:A:220:LEU:HA	1.91	0.41
1:A:437:SER:O	1:A:440:LEU:HG	2.21	0.41
1:A:503:VAL:O	1:A:506:VAL:HG12	2.21	0.41
1:A:260:ILE:HG22	1:A:266:ASN:OD1	2.20	0.41
1:A:260:ILE:C	1:A:266:ASN:OD1	2.59	0.41
1:A:216:TYR:HA	1:A:219:ILE:HG22	2.03	0.41
1:A:256:TYR:CZ	1:A:260:ILE:HD11	2.56	0.41
1:A:334:ILE:O	1:A:346:GLN:NE2	2.53	0.41
1:A:260:ILE:O	1:A:266:ASN:OD1	2.39	0.41
1:A:664:LEU:HD23	1:A:664:LEU:HA	1.93	0.40
1:A:124:GLU:OE2	1:A:191:ARG:NH2	2.50	0.40
1:A:442:ILE:HG23	1:A:471:PHE:CZ	2.56	0.40
1:A:615:ASP:OD1	1:A:616:TYR:N	2.51	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	604/741 (82%)	559 (92%)	40 (7%)	5 (1%)	19 51

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	47	LEU
1	A	99	PRO
1	A	361	PRO
1	A	51	ILE
1	A	647	TRP

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	523/662 (79%)	509 (97%)	14 (3%)	44 71

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	LEU
1	A	83	PHE
1	A	109	LEU

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Mol	Chain	Res	Type
1	A	213	LEU
1	A	269	ASP
1	A	291	ILE
1	A	443	ASN
1	A	471	PHE
1	A	474	ILE
1	A	515	PHE
1	A	532	TYR
1	A	544	LEU
1	A	608	TYR
1	A	657	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 11 ligands modelled in this entry, 11 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, 95^{th} 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	616/741 (83%)	0.55	59 (9%)	8	8	38, 93, 140, 167	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	356	ASN	8.2
1	A	516	ILE	5.2
1	A	249	THR	5.2
1	A	515	PHE	4.9
1	A	39	ASP	4.8
1	A	239	GLU	4.6
1	A	115	LEU	4.3
1	A	248	PHE	4.2
1	A	419	LEU	4.2
1	A	197	LEU	4.2
1	A	390	ASP	4.1
1	A	333	LEU	4.0
1	A	524	PHE	4.0
1	A	463	PRO	3.9
1	A	151	ARG	3.8
1	A	357	TYR	3.7
1	A	399	PHE	3.6
1	A	88	GLU	3.5
1	A	74	TRP	3.5
1	A	241	THR	3.4
1	A	79	PHE	3.3
1	A	514	THR	3.3
1	A	108	TYR	3.3
1	A	527	GLY	3.3
1	A	372	PHE	3.2
1	A	484	TYR	3.0
1	A	366	GLU	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	598	LEU	2.9
1	A	107	TYR	2.9
1	A	525	SER	2.9
1	A	99	PRO	2.8
1	A	367	GLU	2.8
1	A	371	ILE	2.8
1	A	510	GLY	2.8
1	A	64	TYR	2.8
1	A	513	ALA	2.7
1	A	387	GLU	2.7
1	A	253	ALA	2.7
1	A	549	GLN	2.7
1	A	163	LEU	2.6
1	A	148	TRP	2.6
1	A	468	GLU	2.5
1	A	362	LYS	2.5
1	A	150	SER	2.4
1	A	184	PHE	2.4
1	A	116	THR	2.4
1	A	380	ASP	2.3
1	A	349	LYS	2.2
1	A	466	VAL	2.2
1	A	529	TRP	2.2
1	A	257	GLN	2.2
1	A	144	SER	2.1
1	A	375	LEU	2.1
1	A	376	ASP	2.1
1	A	84	LEU	2.1
1	A	66	ILE	2.1
1	A	431	ASN	2.0
1	A	218	ASN	2.0
1	A	346	GLN	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, 95^{th} 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(Å^2)	Q<0.9
3	CA	A	810	1/1	0.33	0.27	118,118,118,118	0
2	BA	A	801	1/1	0.82	0.22	108,108,108,108	1
2	BA	A	806	1/1	0.84	0.11	176,176,176,176	0
2	BA	A	807	1/1	0.90	0.13	163,163,163,163	0
2	BA	A	805	1/1	0.93	0.10	97,97,97,97	0
2	BA	A	802	1/1	0.93	0.29	132,132,132,132	0
2	BA	A	803[A]	1/1	0.94	0.36	72,72,72,72	1
2	BA	A	808	1/1	0.94	0.07	171,171,171,171	0
3	CA	A	809	1/1	0.94	0.04	123,123,123,123	0
2	BA	A	803[B]	1/1	0.94	0.36	72,72,72,72	1
2	BA	A	804	1/1	0.97	0.14	88,88,88,88	0

6.5 Other polymers (i)

There are no such residues in this entry.