Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:06 AM GMT

PDB ID	:	2ZJ7
Title	:	Crystal structure of D157A mutant of Pseudomonas sp. MIS38 lipase
Authors	:	Angkawidjaja, C.; Kuwahara, K.; Kanaya, S.
Deposited on	:	2008-02-29
Resolution	:	2.21 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:

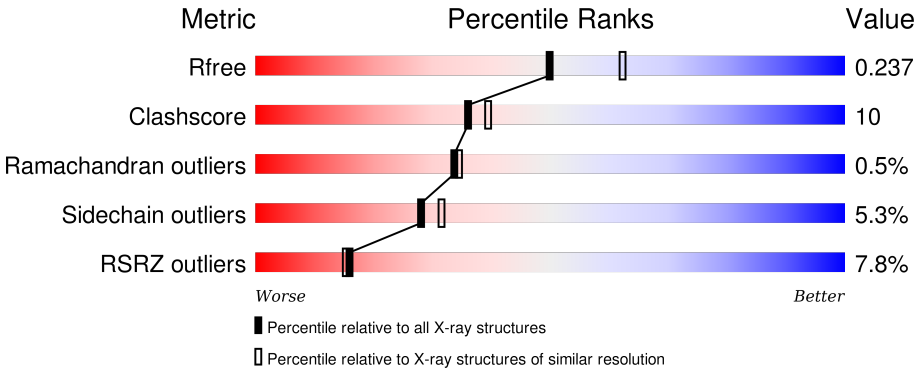
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.21 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	4405 (2.24-2.20)
Clashscore	102246	5146 (2.24-2.20)
Ramachandran outliers	100387	5065 (2.24-2.20)
Sidechain outliers	100360	5066 (2.24-2.20)
RSRZ outliers	91569	4414 (2.24-2.20)

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 <=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	617	8% 81% 16% •

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
3	ZN	A	628	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4997 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 Lipase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	616	Total	C	N	O	S	0	0	0
			4550	2862	769	914	5			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	157	ALA	ASP	ENGINEERED	UNP Q9RBY1

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	10	Total	Ca	0	0
			10	10		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 4 is water.

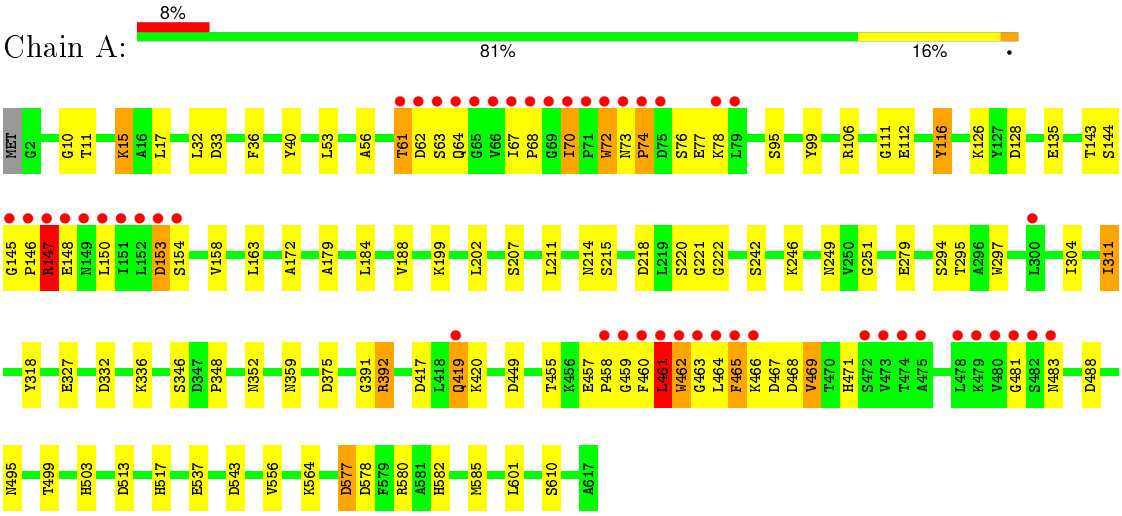
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	435	Total	O	0	0
			435	435		

3 Residue-property plots (i)

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: Lipase

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	48.63Å 84.63Å 87.52Å 90.00° 96.85° 90.00°	Depositor
Resolution (Å)	20.00 – 2.21 40.22 – 2.21	Depositor EDS
% Data completeness (in resolution range)	94.1 (20.00-2.21) 94.1 (40.22-2.21)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.36 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.164 , 0.234 0.181 , 0.237	Depositor DCC
R_{free} test set	1670 reflections (5.27%)	DCC
Wilson B-factor (Å^2)	27.4	Xtriage
Anisotropy	0.098	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 53.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	0 of 33438 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	4997	wwPDB-VP
Average B, all atoms (Å^2)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.99% 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, 0.375 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: ZN, 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	1.03	6/4645 (0.1%)	0.99	16/6311 (0.3%)

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

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	116	TYR	CD1-CE1	5.76	1.48	1.39
1	A	207	SER	CB-OG	-5.67	1.34	1.42
1	A	99	TYR	CD2-CE2	-5.55	1.31	1.39
1	A	279	GLU	CG-CD	5.55	1.60	1.51
1	A	332	ASP	CB-CG	-5.38	1.40	1.51
1	A	77	GLU	CG-CD	5.03	1.59	1.51

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	392	ARG	NE-CZ-NH1	-11.64	114.48	120.30
1	A	392	ARG	NE-CZ-NH2	9.44	125.02	120.30
1	A	419	GLN	CB-CA-C	-7.42	95.56	110.40
1	A	222	GLY	N-CA-C	-6.79	96.13	113.10
1	A	221	GLY	N-CA-C	6.43	129.16	113.10
1	A	336	LYS	CD-CE-NZ	-6.10	97.67	111.70
1	A	221	GLY	C-N-CA	5.82	134.53	122.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	464	LEU	CA-CB-CG	5.79	128.61	115.30
1	A	33	ASP	CB-CG-OD1	5.65	123.39	118.30
1	A	128	ASP	CB-CG-OD1	5.61	123.35	118.30
1	A	417	ASP	CB-CG-OD1	5.46	123.22	118.30
1	A	449	ASP	CB-CG-OD1	-5.24	113.58	118.30
1	A	15	LYS	CD-CE-NZ	-5.20	99.73	111.70
1	A	543	ASP	CB-CG-OD1	5.20	122.98	118.30
1	A	221	GLY	CA-C-N	5.10	126.40	116.20
1	A	128	ASP	CB-CG-OD2	-5.03	113.78	118.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	143	THR	Peptide
1	A	145	GLY	Peptide
1	A	461	LEU	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	4550	0	4340	85	1
2	A	10	0	0	0	0
3	A	2	0	0	0	2
4	A	435	0	0	11	1
All	All	4997	0	4340	85	3

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 10.

All (85) 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:146:PRO:HB2	1:A:148:GLU:HB2	1.25	1.14

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63:SER:HB2	1:A:64:GLN:HA	1.13	1.07
1:A:146:PRO:CB	1:A:148:GLU:HB2	1.85	1.05
1:A:95:SER:HB3	4:A:840:HOH:O	1.58	1.02
1:A:462:TRP:HB3	1:A:463:GLY:HA3	1.45	0.98
1:A:147:ARG:HG3	1:A:147:ARG:HH11	1.30	0.96
1:A:63:SER:CB	1:A:64:GLN:HA	1.97	0.94
1:A:462:TRP:HB3	1:A:463:GLY:CA	1.98	0.93
1:A:72:TRP:HA	1:A:73:ASN:HB2	1.57	0.86
1:A:458:PRO:N	1:A:459:GLY:HA2	1.94	0.83
1:A:70:ILE:HG22	4:A:990:HOH:O	1.78	0.82
1:A:457:GLU:C	1:A:459:GLY:HA2	2.00	0.82
1:A:153:ASP:HA	4:A:820:HOH:O	1.79	0.82
1:A:146:PRO:HB2	1:A:148:GLU:CB	2.10	0.79
1:A:214:ASN:HD21	1:A:249:ASN:HD21	1.32	0.78
1:A:462:TRP:CB	1:A:463:GLY:HA3	2.14	0.77
1:A:147:ARG:CG	1:A:147:ARG:HH11	1.99	0.73
1:A:10:GLY:HA3	4:A:654:HOH:O	1.87	0.73
1:A:294:SER:HB3	1:A:297:TRP:HB2	1.75	0.69
1:A:36:PHE:HB2	1:A:70:ILE:HG23	1.76	0.68
1:A:327:GLU:OE1	1:A:392:ARG:NH1	2.22	0.68
1:A:146:PRO:HB3	1:A:148:GLU:HB2	1.75	0.67
1:A:460:PHE:N	1:A:461:LEU:HA	2.12	0.63
1:A:32:LEU:HD11	1:A:56:ALA:HB1	1.81	0.63
1:A:462:TRP:CE3	1:A:462:TRP:HA	2.35	0.62
1:A:147:ARG:O	1:A:148:GLU:CG	2.49	0.61
1:A:150:LEU:HB2	4:A:767:HOH:O	2.03	0.59
1:A:64:GLN:HE22	1:A:76:SER:HB2	1.68	0.59
1:A:577:ASP:OD2	1:A:582:HIS:CE1	2.56	0.58
1:A:63:SER:HB2	1:A:64:GLN:CA	2.08	0.58
1:A:146:PRO:C	1:A:148:GLU:N	2.57	0.57
1:A:147:ARG:O	1:A:148:GLU:HG2	2.04	0.57
1:A:211:LEU:HD23	1:A:211:LEU:C	2.24	0.56
1:A:147:ARG:C	4:A:767:HOH:O	2.42	0.56
1:A:495:ASN:HD22	1:A:513:ASP:H	1.52	0.56
1:A:147:ARG:HG3	1:A:147:ARG:NH1	2.11	0.55
1:A:462:TRP:HE3	1:A:462:TRP:HA	1.71	0.55
1:A:172:ALA:O	1:A:215:SER:HB2	2.08	0.54
1:A:577:ASP:OD2	1:A:582:HIS:HE1	1.91	0.54
1:A:251:GLY:HA2	1:A:318:TYR:HE1	1.72	0.54
1:A:111:GLY:HA3	1:A:116:TYR:O	2.07	0.53
1:A:147:ARG:C	1:A:148:GLU:CG	2.77	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:455:THR:OG1	1:A:471:HIS:HE1	1.92	0.53
1:A:64:GLN:HE22	1:A:76:SER:CB	2.22	0.53
1:A:147:ARG:CG	1:A:147:ARG:NH1	2.64	0.52
1:A:147:ARG:C	1:A:148:GLU:HG3	2.30	0.52
1:A:462:TRP:HB3	1:A:463:GLY:C	2.29	0.52
1:A:73:ASN:N	1:A:74:PRO:HD3	2.24	0.51
1:A:184:LEU:O	1:A:188:VAL:HG23	2.10	0.51
1:A:147:ARG:HB3	4:A:711:HOH:O	2.10	0.51
1:A:461:LEU:HB3	1:A:462:TRP:CE3	2.46	0.50
1:A:466:LYS:HD2	4:A:1061:HOH:O	2.12	0.49
1:A:11:THR:O	1:A:15:LYS:HG3	2.12	0.49
1:A:391:GLY:O	1:A:392:ARG:HB2	2.12	0.48
1:A:153:ASP:HB3	1:A:158:VAL:HG23	1.94	0.48
1:A:375:ASP:OD1	1:A:392:ARG:HD3	2.13	0.48
1:A:17:LEU:HD22	1:A:202:LEU:HD13	1.96	0.47
1:A:457:GLU:HG2	1:A:469:VAL:CG1	2.45	0.46
1:A:126:LYS:HB3	1:A:135:GLU:HB2	1.98	0.46
1:A:455:THR:O	1:A:468:ASP:HA	2.17	0.45
1:A:61:THR:N	1:A:147:ARG:HE	2.15	0.45
1:A:462:TRP:CD1	1:A:465:PHE:O	2.70	0.45
1:A:67:ILE:O	1:A:70:ILE:HD12	2.17	0.45
1:A:220:SER:OG	1:A:246:LYS:HE3	2.16	0.45
1:A:578:ASP:OD2	1:A:580:ARG:NE	2.50	0.45
1:A:72:TRP:HA	1:A:73:ASN:CB	2.37	0.45
1:A:61:THR:CA	1:A:147:ARG:HE	2.30	0.44
1:A:153:ASP:CB	1:A:158:VAL:HG23	2.47	0.44
1:A:419:GLN:O	1:A:420:LYS:HG2	2.18	0.44
1:A:503:HIS:HB2	4:A:633:HOH:O	2.17	0.44
1:A:40:TYR:CE1	1:A:311:ILE:HD11	2.53	0.44
1:A:53:LEU:HB2	1:A:70:ILE:HG12	2.00	0.43
1:A:64:GLN:NE2	1:A:76:SER:HB2	2.31	0.43
1:A:106:ARG:HD3	4:A:882:HOH:O	2.18	0.43
1:A:218:ASP:OD1	1:A:242:SER:HB2	2.18	0.43
1:A:67:ILE:HA	1:A:68:PRO:HD3	1.84	0.43
1:A:556:VAL:HB	1:A:601:LEU:HD23	2.01	0.43
1:A:499:THR:HA	1:A:517:HIS:O	2.19	0.42
1:A:112:GLU:HG3	1:A:179:ALA:O	2.20	0.42
1:A:459:GLY:HA3	1:A:460:PHE:C	2.40	0.41
1:A:163:LEU:HD11	1:A:304:ILE:HD13	2.02	0.41
1:A:147:ARG:O	1:A:148:GLU:HG3	2.21	0.41
1:A:457:GLU:OE1	1:A:457:GLU:HA	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:TRP:HB3	4:A:748:HOH:O	2.21	0.40
1:A:346:SER:HB2	1:A:348:PRO:HD2	2.02	0.40

All (3) 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:537:GLU:OE2	3:A:628:ZN:ZN[2_656]	1.21	0.99
3:A:628:ZN:ZN	3:A:629:ZN:ZN[2_646]	1.38	0.82
4:A:860:HOH:O	4:A:1054:HOH:O[2_545]	2.16	0.04

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	614/617 (100%)	594 (97%)	17 (3%)	3 (0%)	34 34

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	147	ARG
1	A	74	PRO
1	A	481	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	469/470 (100%)	444 (95%)	25 (5%)	28 31

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	61	THR
1	A	62	ASP
1	A	70	ILE
1	A	72	TRP
1	A	78	LYS
1	A	144	SER
1	A	147	ARG
1	A	153	ASP
1	A	154	SER
1	A	199	LYS
1	A	295	THR
1	A	311	ILE
1	A	352	ASN
1	A	359	ASN
1	A	461	LEU
1	A	462	TRP
1	A	465	PHE
1	A	467	ASP
1	A	469	VAL
1	A	483	ASN
1	A	488	ASP
1	A	564	LYS
1	A	577	ASP
1	A	585	MET
1	A	610	SER

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	64	GLN
1	A	249	ASN
1	A	298	ASN
1	A	352	ASN
1	A	471	HIS
1	A	495	ASN
1	A	553	GLN
1	A	582	HIS

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

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	616/617 (99%)	0.34	48 (7%)	16	15	13, 27, 78, 111	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	462	TRP	14.4
1	A	461	LEU	12.8
1	A	72	TRP	12.5
1	A	464	LEU	11.6
1	A	460	PHE	10.5
1	A	465	PHE	10.4
1	A	152	LEU	9.5
1	A	68	PRO	9.4
1	A	459	GLY	8.8
1	A	70	ILE	8.6
1	A	151	ILE	8.6
1	A	149	ASN	8.5
1	A	66	VAL	8.3
1	A	463	GLY	8.1
1	A	71	PRO	6.6
1	A	64	GLN	6.5
1	A	73	ASN	6.4
1	A	153	ASP	6.3
1	A	69	GLY	6.2
1	A	150	LEU	5.8
1	A	63	SER	5.7
1	A	74	PRO	5.5
1	A	145	GLY	5.1
1	A	466	LYS	5.0
1	A	458	PRO	5.0
1	A	146	PRO	4.8
1	A	474	THR	4.5

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Mol	Chain	Res	Type	RSRZ
1	A	65	GLY	4.2
1	A	67	ILE	4.2
1	A	147	ARG	4.2
1	A	75	ASP	4.0
1	A	483	ASN	3.8
1	A	62	ASP	3.8
1	A	61	THR	3.6
1	A	479	LYS	3.3
1	A	475	ALA	3.2
1	A	148	GLU	2.9
1	A	79	LEU	2.9
1	A	473	VAL	2.8
1	A	478	LEU	2.8
1	A	154	SER	2.7
1	A	482	SER	2.7
1	A	481	GLY	2.7
1	A	300	LEU	2.5
1	A	78	LYS	2.5
1	A	480	VAL	2.3
1	A	419	GLN	2.1
1	A	472	SER	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, 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	LLDF	B-factors(Å^2)	Q<0.9
3	ZN	A	628	1/1	1.00	0.11	-0.21	150,150,150,150	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
2	CA	A	627	1/1	0.92	0.09	-0.94	29,29,29,29	0
2	CA	A	621	1/1	0.99	0.08	-0.94	26,26,26,26	0
2	CA	A	618	1/1	0.99	0.08	-1.46	19,19,19,19	0
2	CA	A	624	1/1	0.99	0.08	-1.59	22,22,22,22	0
2	CA	A	626	1/1	0.99	0.08	-1.69	27,27,27,27	0
2	CA	A	625	1/1	0.98	0.08	-1.78	28,28,28,28	0
2	CA	A	619	1/1	1.00	0.07	-1.79	18,18,18,18	0
2	CA	A	620	1/1	0.99	0.06	-2.49	25,25,25,25	0
2	CA	A	622	1/1	1.00	0.07	-2.50	27,27,27,27	0
2	CA	A	623	1/1	0.99	0.04	-3.19	27,27,27,27	0
3	ZN	A	629	1/1	0.99	0.06	-4.05	31,31,31,31	0

6.5 Other polymers

There are no such residues in this entry.